



REMEDIAL ACTION QUARTERLY MONITORING REPORT

**FIRST QUARTER – 2004
(3 of 120)**

**SKINNER LANDFILL SITE
BUTLER COUNTY
WEST CHESTER, OHIO**

Prepared for:

Skinner Landfill Work Group
c/o Ben Baker
The Dow Chemical Company
Michigan Operations
47 Building
Midland, MI 48667

Prepared by:

Earth Tech, Inc.
2373 Progress Drive
Hebron, KY 41048

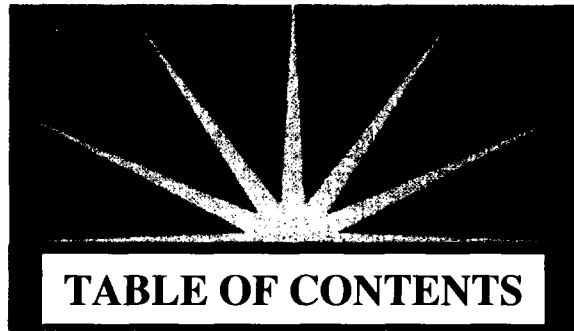


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**LIST OF
ACRONYMS**



LIST OF ACRONYMS

| | |
|--------|--|
| AMP | Air Monitoring Plan |
| AOC | Administrative Order on Consent |
| ARAR | Applicable or Relevant and Appropriate Requirements |
| BMR | Baseline Monitor Report |
| BCDES | Butler County Department of Environmental Services |
| bgs | Below Ground Surface |
| BZ | Breathing Zone |
| CD&D | Construction Debris and Demolition Waste |
| CERCLA | Comprehensive Environmental Response, Compensation and Liability Act |
| CGI | Combustible Gas Indicator |
| CHSD | Corporate Health and Safety Director |
| CIP | Construction Implementation Plan |
| CLP | Contract Laboratory Program |
| cm/sec | Centimeters Per Second |
| CO | Carbon Monoxide |
| CP | Contingency Plan |
| CQA | Construction Quality Assurance |
| CQAC | Construction Quality Assurance Consultant |
| CRZ | Contamination Reduction Zone |
| CRQL | Contract Required Quantitation Limit |
| CSDI | Contaminated Soils Design Investigation |
| CY | Cubic Yard |
| CZ | Control Zone |
| DSW | Division of Surface Water (OEPA) |
| DSR | Division Safety Representative |
| EPA | Environmental Protection Agency |
| EZ | Exclusion Zone |
| FID | Flame Ionization Detector |
| FML | Flexible Membrane Liner (low density polyethylene) |
| FSP | Field Sampling Plan |
| FTB | Film Tearing Bond |
| ft | Feet |
| ft/sec | Feet Per Second |
| GCL | Geosynthetic Clay Layer |
| GCAL | Gulf Coast Analytical Laboratories Inc. |
| GIS | Groundwater Interceptor System |
| gpd | Gallons Per Day |
| gpm | Gallons Per Minute |
| GWDI | Groundwater Design Investigation |
| HAP | Hazardous Air Pollutant |
| HASP | Health and Safety Plan |
| HDPE | High-Density Polyethylene |
| HSM | Health and Safety Manager |
| IDLH | Immediately Dangerous to Life or Health |

| | |
|-----------------|---|
| IRM | Interim Remedial Measures |
| kg/d | Kilograms Per Day |
| lb/day | Pounds Per Day |
| LEL | Lower Explosion Limit |
| LF | Lineal Feet |
| LLDPE | Linear Low-Density Polyethylene |
| μ | Micron |
| μg/l | Microgram per Liter |
| MSL | Mean Sea Level |
| NIOSH | National Institute for Occupational Safety and Health |
| NO _x | Oxides of Nitrogen |
| NWI | National Wetland Inventory |
| O ₃ | Ozone |
| OAC | Ohio Administrative Code |
| ODNR | Ohio Department of Natural Resources |
| OEPA | Ohio Environmental Protection Agency |
| ORC | Ohio Revised Code |
| OSHA | Occupational Safety and Health Administration |
| PEL | Permissible Exposure Limit |
| PID | Photoionization Detector |
| PLC | Programmable Logic Controller |
| PM-10 | Particulate Matter less than 10 microns |
| PRP | Potentially Responsible Party |
| PPE | Personal Protective Equipment |
| psi | Pounds Per Square Inch |
| PQL | Practical Quantitation Limit |
| QAPP | Quality Assurance Project Plan |
| QA | Quality Assurance |
| QC | Quality Control |
| RCRA | Resource Conservation and Recovery Act |
| RA | Remedial Action |
| RD | Remedial Design |
| RHSS | Regional Health & Safety Specialist |
| RI/FS | Remedial Investigation/Feasibility Study |
| ROD | Record of Decision |
| RPM | Remedial Project Manager (USEPA) |
| RPO | Resident Project Observer |
| SI | Site Inspection |
| SF | Square Feet |
| SLWG | Skinner Landfill Work Group |
| SO ₂ | Sulfur Dioxide |
| SOP | Standard Operating Procedure |
| SOW | Statement of Work |
| SPCC | Spill Prevention Control and Counter Measure Plan |
| SSO | Site Safety Officer |
| SVE | Soil Vapor Extraction |
| SVOC | Semi-Volatile Organic Compound |
| SZ | Support Zone |

| | |
|-------|---|
| TAL | Target Analyte List |
| TCL | Target Compound List |
| TDH | Total Dynamic Head |
| TLV | Threshold Limit Values |
| TSS | Total Suspended Solids |
| TWA | Time Weighted Average |
| USACE | United States Army Corps of Engineers |
| USEPA | United States Environmental Protection Agency |
| USFWS | United States Fish and Wildlife Services |
| USGS | United States Geological Survey |
| VOC | Volatile Organic Compound |
| yr | Year |
| WBG | Wet Bulb Globe Temperature |
| WZ | Work Zone |

1.0 INTRODUCTION

1.1 GENERAL INFORMATION

This quarterly monitoring report was prepared for the Skinner Landfill Superfund Site located in West Chester, Butler County, Ohio in accordance with the Operation and Maintenance - Long-Term Performance Plan (O&M-LTP Plan) dated August 2003. The O&M-LTP Plan was prepared to meet the requirements of the Record of Decision (ROD) dated June 4, 1993, the Statement of Work (SOW) dated April 6, 1994, the 100% Final Remedial Design dated June 21, 1996 and the Consent Decree dated April 7, 2001.

The remedial action (RA) post-construction O&M monitoring period began with the third quarter of 2003 and extends for a period of 30 years. This report documents the results of groundwater and surface water monitoring conducted during the first quarter of 2004, which is the 3rd of 120 sampling events to be conducted during the 30-year monitoring period.

1.2 SITE LOCATION AND DESCRIPTION

Skinner Landfill is located approximately 15 miles north of Cincinnati, Ohio near West Chester, Butler County, Ohio in Township 3, Section 22, Range 2. The site is located along Cincinnati-Dayton Road, as shown in Figure 1. The site is bordered on the south by the East Fork of Mill Creek, on the north by wooded land, on the east by a Norfolk Southern Railway Company right-of-way, and on the west by a gravel driveway.

The site is located in a highly dissected area that slopes from a till-mantled-bedrock upland to a broad, flat-bottomed valley that is occupied by the main branch of Mill Creek. Elevations on the site range from a high of nearly 300 feet above mean sea level (MSL) in the northeast, to a low of 645 feet above MSL near the confluence of Skinner Creek and East Fork of Mill Creek. Both Skinner Creek and the East Fork of Mill Creek are small, intermittent shallow streams. Both of these streams flow to the southwest from the site toward the main branch of Mill Creek.

In general, the site is underlain by relatively thin glacial drift over inter-bedded shale and limestone of Ordovician age. The composition of the glacial drift ranges from intermixed silt, sand and gravel, to silty sandy clays with a thickness ranging from zero to over forty feet. The sand and gravel deposits comprise the hills and ridges and are encountered near the surface of the central portion of the site. The silts and clays usually occur as lenses in the sands and gravel or directly overlie bedrock.

1.3 SITE HISTORY AND BACKGROUND

The property was originally developed as a sand and gravel mining operation and was subsequently used as a landfill from 1934 to 1990. According to USEPA studies, materials deposited at the site include demolition debris, household refuse and a wide variety of chemical wastes. The waste disposal areas include a now buried former waste lagoon near the center of the site and a landfill. According to USEPA studies, the buried lagoon was used for the disposal of paint wastes, ink wastes, creosote, pesticides, and other chemical wastes. The landfill area, located north and northeast of the buried lagoon, received predominantly demolition and landscaping debris.

In 1976, the Ohio EPA (OEPA) initiated an investigation of the site. In 1982, the site was placed on the National Priority List by the USEPA based on information obtained during a limited investigation of the

site. A Phase II Remedial Investigation was conducted from 1989 to 1991 and involved further investigation of groundwater, surface water, soils and sediments. Both a Baseline Risk Assessment and Feasibility Study (FS) were completed in 1992.

The Phase II Remedial Investigation revealed that the most contaminated media at the site is the soil in the buried waste lagoon. Migration of the landfill constituents has been limited, and the Phase II Remedial Investigation concluded that there had been no off-site migration of landfill constituents via groundwater flow.

In the Record of Decision (ROD), dated June 4, 1993, the USEPA selected a remedy for the site consisting of multi-media capping of the landfill and the buried waste lagoon, and collection and treatment of the groundwater. The ROD also required an investigation to determine the feasibility for soil vapor extraction (SVE) in the granular soil adjacent to the buried lagoon.

The Remedial Design (RD) Investigation performed in 1994 was implemented to collect data required to assess the feasibility of the SVE and to design the multi-media cap and the groundwater extraction/treatment systems. The Remedial Design was submitted to USEPA on June 21, 1996 outlining the cover design and groundwater interception system design. Based on the RD investigation, the installation of an SVE system was determined to be unfeasible.

Construction of a groundwater interception system (GIS) and engineered landfill cover system began in April 2001 and was substantially completed in September 2001. The USEPA conducted the pre-final construction inspection on September 27, 2001, the final construction inspection on March 27, 2003 and the second 5-Year Review on January 22, 2004.

2.0 SAMPLING METHODS

This quarterly monitoring event was conducted in general accordance with the following documents shown with the date of the USEPA-approved final version:

- Operation and Maintenance - Long-Term Performance Plan (O&M-LTP Plan) dated August 2003, and
- RA Health and Safety Plan, Final February 2001.

There were no deviations from these work plans.

3.0 RESULTS

3.1 GROUNDWATER LEVELS

The groundwater elevation data obtained from the monitor wells, piezometers and selected gas probes is presented on Table 1 with the corresponding potentiometric surface map provided in Appendix A. The groundwater flow direction and gradient remained relatively unchanged when compared to the previous quarterly monitoring report period. Groundwater flow direction is to the south-southeast directly toward the East Fork of Mill Creek with an average hydraulic gradient of 0.12 ft/ft. The groundwater gradient has remained relatively unchanged when compared to the average hydraulic gradient of 0.13 ft/ft documented in the Remedial Action Baseline Monitoring Report dated March 2005.

3.2 GROUNDWATER-WASTE MONITORING

Results of the piezometer groundwater levels used to monitor the groundwater levels relative to bottom of waste are provided on Table 2. Based on measured water levels, groundwater has been lowered below the waste elevation during this monitor event at piezometers P-11 and P-12, which are the two piezometers furthest from Duck Pond. The groundwater level remains above the bottom of waste at piezometers P-9 and P-10. The depth to water measurements for piezometers P-9, P-10 and P-11 were obtained with a smaller diameter water level indicator, as opposed to a groundwater interface probe, due to a pinching of the well casings that reduced the diameter of the piezometers.

3.3 GROUNDWATER ANALYTICAL RESULTS

A summary of target compound list (TCL) and target analyte list (TAL) parameter concentrations encountered above the contract required detection limit and revised modified trigger level is provided on Table 3. A summary of the laboratory analytical results have been presented on a per well basis in Appendix B to assist in identifying temporal detection patterns. A report of each data set reduction, validation and assessment procedure conducted on an analytical-set basis in accordance with the O&M-LTP Plan quality assurance project plan (QAPP) is included in Appendix C.

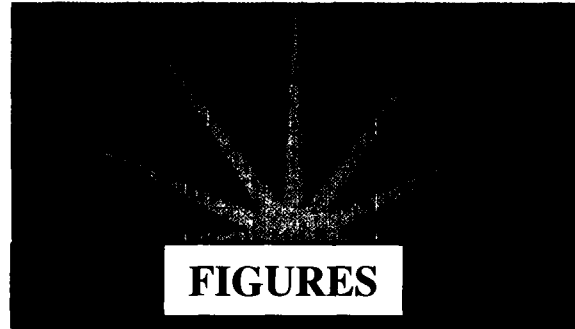
In general, target compound list volatiles, semi-volatiles, pesticides and PCBs were not detected in groundwater above the CRQL.

Two of the sixteen TAL parameters, with a revised modified trigger level, were detected above the CRQL. Detections of iron (present in two groundwater monitoring wells) and barium (present in one groundwater monitoring well) were detected above the CRQL, but below the revised modified trigger levels.

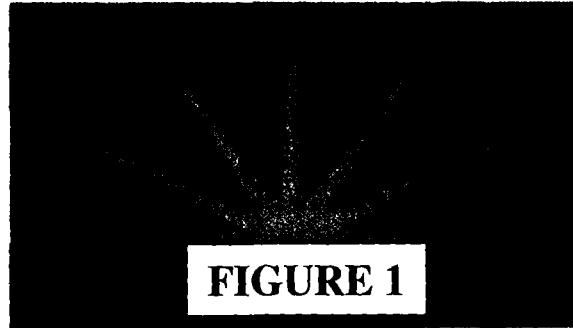
3.4 SURFACE WATER ANALYTICAL RESULTS

Surface water analyzed consisted of sampling surface runoff from the site and surface water directly from the East Fork of Mill Creek. A summary of TCL and TAL parameter concentrations encountered above the contract required detection limit and revised modified trigger level is provided on Table 4. A summary of surface water laboratory analytical results is presented in Appendix B. The summary tables are presented on a sample location basis. The validated laboratory analytical data is provided in Appendix C.

In general, target compound list volatiles, semi-volatiles, pesticides and PCBs were not detected in surface water above the CRQL. None of the sixteen TAL parameters, with a revised modified trigger level, were detected above the CRQL.

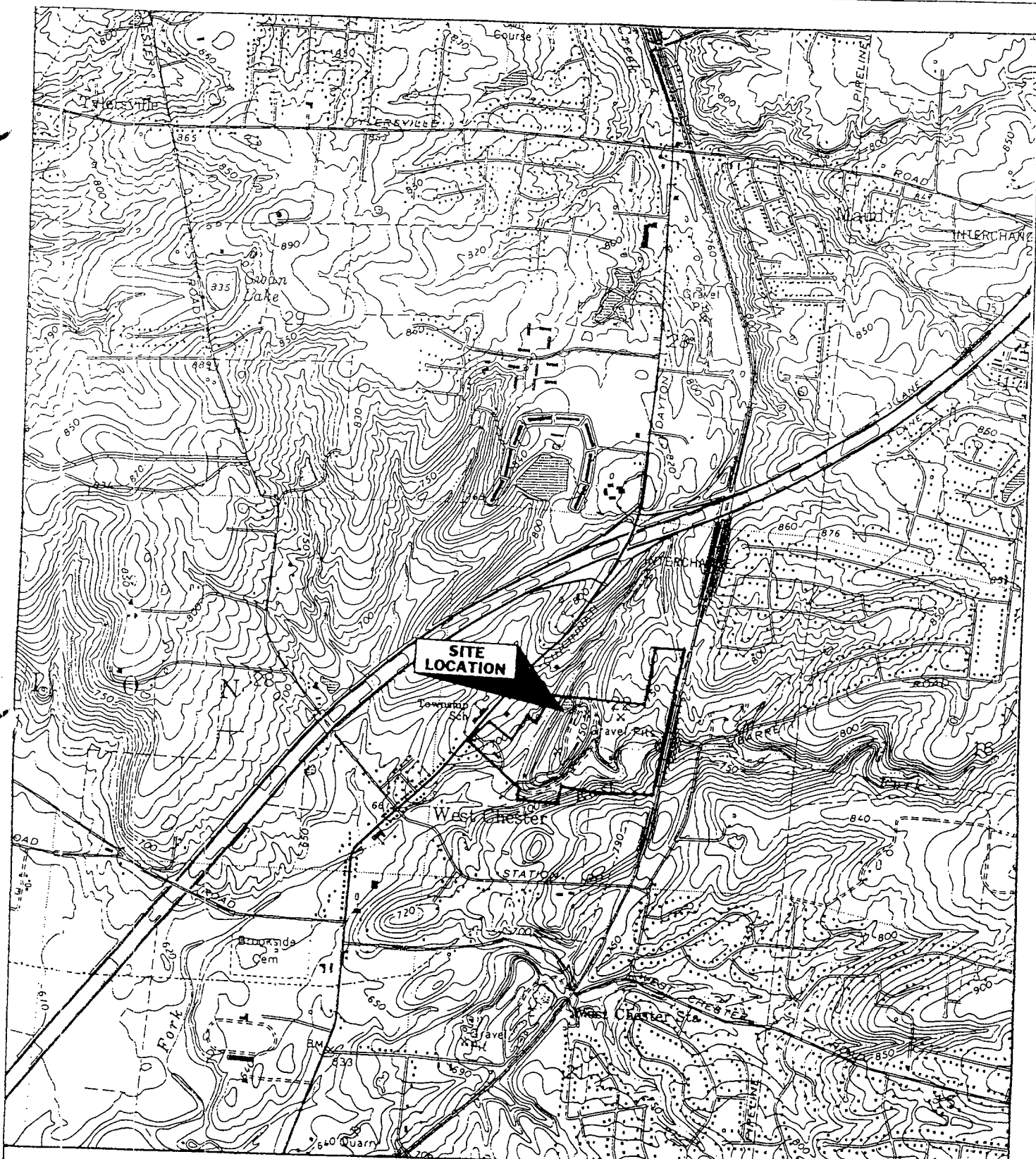


FIGURES



SITE VICINITY MAP

FIGURE 1



Base taken from USGS Glendale, Ohio
7.5' Topographic Quadrangle, photorevised 1987



EARTH TECH



SKINNER LANDFILL
SITE VICINITY MAP
BUTLER COUNTY, OHIO



TABLE 1

**GROUNDWATER
ELEVATIONS**

TABLE 1

TABLE 1
Groundwater Elevation Summary
Skinner Landfill
West Chester, Ohio

| Well Type | Location | Well Use | Ground Surface Elevation (MSL-feet) | Top of Casing Elevation (MSL-feet) | March 2004 | |
|---------------------------------|----------|----------|--|---------------------------------------|---|-------------------------------------|
| | | | | | Depth to Water (feet from top of casing) | Groundwater Elevation (MSL-feet) |
| Piezometers | P-1 | G | 685.42 | 687.65 | 9.75 | 677.90 |
| | P-2 | G | 688.54 | 690.42 | 13.21 | 677.21 |
| | P-3R | G | 691.83 | 693.69 | 25.43 | 668.26 |
| | P-4 | G | 700.32 | 702.63 | 6.68 | 695.95 |
| | P-5 | G | 708.20 | 710.65 | 14.33 | 696.32 |
| | P-6 | G | 707.45 | 710.59 | 12.45 | 698.14 |
| | P-7 | G | 719.08 | 721.83 | Dry | Dry |
| | P-8 | G | 747.70 | 749.91 | 29.73 | 720.18 |
| | P-9 | G | 760.68 | 763.90 | 20.75 | 743.15 |
| | P-10 | G | 761.34 | 764.16 | 26.10 | 738.06 |
| | P-11 | G | 760.34 | 762.76 | 27.40 | 735.36 |
| | P-12 | G | 743.50 | 746.17 | 40.65 | 705.52 |
| Groundwater Monitoring Wells | GW-06R | S | 683.89 | 685.91 | 9.94 | 675.97 |
| | GW-07R | S | 683.46 | 683.06 | 4.31 | 678.75 |
| | GW-24 | G | 693.32 | 695.21 | 17.65 | 677.56 |
| | GW-26 | G | 696.61 | 698.28 | 29.82 | 668.46 |
| | GW-30 | G | 675.63 | 677.62 | 10.18 | 667.44 |
| | GW-58 | S | 684.03 | 686.53 | 14.13 | 672.40 |
| | GW-59 | S | 684.35 | 687.38 | 6.97 | 680.41 |
| | GW-60 | S | 689.12 | 692.38 | 13.08 | 679.30 |
| | GW-61 | S | 687.38 | 690.86 | 12.88 | 677.98 |
| | GW-62A | S | 690.19 | 692.38 | 29.81 | 662.57 |
| | GW-62B | S | 690.57 | 693.13 | 12.48 | 680.65 |
| | GW-63 | S | 698.87 | 702.50 | 10.23 | 692.27 |
| | GW-64 | S | 700.45 | 703.88 | 13.17 | 690.71 |
| | GW-65 | S | 703.83 | 706.88 | 13.64 | 693.24 |
| | GW-66 | G | 686.82 | 689.41 | 8.67 | 680.74 |
| Gas Probes | GP-6 | G | 772.18 | 774.65 | 14.91 | 759.74 |
| | GP-7 | G | 749.83 | 752.65 | 8.62 | 744.03 |

Notes:

MSL - Mean Sea Level

G - Gauging

S - Sampling and Gauging

— No Gauging Data Available (well constricted)



**GROUNDWATER/WASTE
ELEVATIONS**

TABLE 2

TABLE 2
Groundwater-Waste Monitoring Summary
Skinner Landfill
West Chester, Ohio

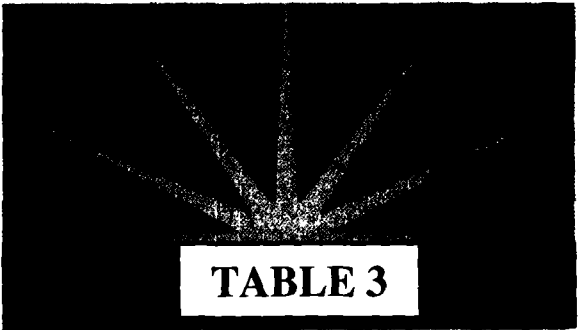
| Piezometer | Depth to Waste (feet) | Bottom of Waste Elevation (MSL-feet) | Baseline Water Elevation (June 2001) (feet) | Water Elevation (May 2003) (feet) | Water Elevation (August 2003) (feet) | Water Elevation (November 2003) (feet) | Water Elevation (March 2004) (feet) |
|------------|--------------------------|--|--|---|--|--|---|
| P-9 | 25 | 737 | 745.00 | 744.75 | - | - | 743.15 |
| P-10 | 30 | 734 | 744.50 | 744.71 | 738.46 | 739.16 | 738.06 |
| P-11 | 17 | 745 | 744.30 | 734.82 | 735.74 | 734.83 | 735.36 |
| P-12 | 35 | 707 | 713.50 | 705.83 | 705.42 | 705.39 | 705.52 |

Notes:

Waste elevations determined during piezometer installation on June 28 and 29, 2001.

Shaded cells indicate water level elevations below the elevation of waste.

- No Gauging Data Available (well constricted)



**GROUNDWATER RESULTS
SUMMARY**

Table 3

Groundwater Summary

Skinner Landfill
West Chester, Ohio
First Quarter 2004

| Sample ID | VOCs | SVOCs | Dissolved Metals** | Pesticides/PCBs |
|-----------|------|-------|--------------------|-----------------|
| GW-06R | - | - | <i>barium</i> | - |
| GW-07R | - | - | - | - |
| GW-58 | - | - | <i>iron</i> | - |
| GW-59 | - | - | - | - |
| GW-60 | - | - | - | - |
| GW-61 | - | - | <i>iron</i> | - |
| GW-62A | - | - | - | - |
| GW-62B | * | * | * | * |
| GW-63 | - | - | - | - |
| GW-64 | - | - | - | - |
| GW-65 | - | * | * | * |

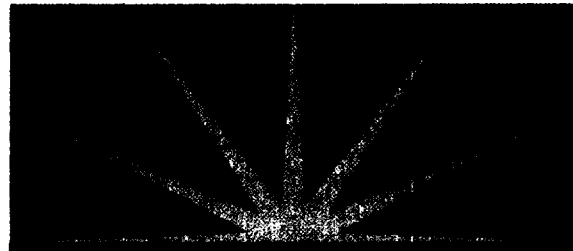
- all parameters below report limits

italic - above Contract Required Quantitation Levels (CRQL's)

bold - above trigger level

* - Insufficient sample volume.

** - Dissolved metals for analytes that have a corresponding trigger level.



| |
|---------|
| TABLE 4 |
|---------|

**SURFACE WATER
RESULTS SUMMARY**

Table 4

Surface Water Summary

Skinner Landfill
West Chester, Ohio
First Quarter 2004

| Sample ID | VOCs | SVOCs | Dissolved Metals** | Pesticides/PCBs |
|-----------|------|-------|--------------------|-----------------|
| SW-50 | - | - | - | - |
| SW-51 | - | - | - | - |
| SW-52 | - | - | - | - |
| SWD-1 | * | * | * | * |
| SWD-2 | * | * | * | * |
| SWD-3 | - | - | - | - |

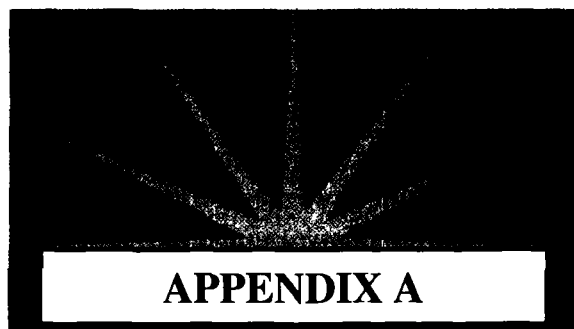
- all parameters below report limits

italic - above Contract Required Quantitation Levels (CRQL's)

bold - above trigger level

* - Insufficient sample volume.

** - Dissolved metals for analytes that have a corresponding trigger level.



APPENDIX A

POTENTIONMETRIC SURFACE MAP

SDMS US EPA Region V

Imagery Insert Form



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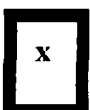
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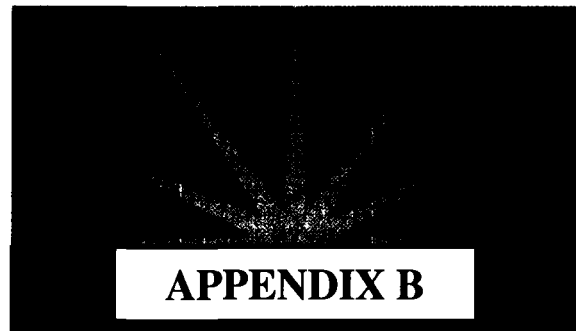
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APPENDIX A – POTENTIOMETRIC SURFACE MAP



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SUMMARY OF ANALYTICAL RESULTS

APPENDIX B

**Skinner Landfill
West Chester, Ohio
Groundwater Analysis Summary Table for Monitoring Well GW-06R**

| Sampling Event (All Results Expressed in Units of µg/l) | | | | | | | | | |
|---|------------------|-------------|-------------|--------|-------------------|-------------|----------|---------------|-------|
| Compound | Baseline Results | | | | Quarterly Results | | | TRIGGER LEVEL | CRQL |
| | September-02 | December-02 | February-03 | May-03 | August-03 | November-03 | March-04 | | |
| Inorganics - Metals (Dissolved) ¹³ | | | | | | | | | |
| Aluminum | — | — | — | — | — | 25.8 | 25.8 | | 200 |
| Antimony | 13.3 B | 6.0 B | 3.0 | 7.0 | 3.7 | 3.7 | 3.7 | 60 | 60 |
| Arsenic | 8.1 B | 3.6 U | 3.6 | 2.9 | 2.9 | 2.9 | 2.9 | 20 | 10 |
| Barium | 266 | 254 | 256 | 224 | 309 J | 294 | 266 | 1,000 | 200 |
| Beryllium | 0.1 U | 0.1 U | 0.1 | 0.1 | 0.1 | 0.1 | 0.1 | 5 | 5 |
| Cadmium | 0.2 U | 0.2 U | 0.2 UJ | 0.2 | 0.2 | 0.2 | 0.2 | 5 | 5 |
| Calcium | — | — | — | — | — | 189,000 | 189,000 | | 5,000 |
| Chromium | 1.1 B | 3.5 B | 2.1 | 2.6 | 2.6 | 0.8 | 1.2 | 11 | 10 |
| Cobalt | — | — | — | — | — | 0.4 | 0.4 | | 50 |
| Copper | 1.0 U | 1.0 U | 1.0 UJ | 1.9 | 1.3 | 1.7 | 1.2 | 25 | 25 |
| Iron | 92.2 B | 4.9 U | 79.2 | 14.1 | 14.1 | 14.1 | 22.0 | 7,000 | 100 |
| Lead | 1.3 U | 1.3 UJ | 1.3 R | 1.5 UJ | 1.5 | 1.5 | 1.5 | 4.2 | 3 |
| Magnesium | — | — | — | — | — | 30,500 | 30,000 | | 5,000 |
| Manganese | — | — | — | — | — | 77.0 | 69.5 J | | 15 |
| Mercury | 0.1 U | 0.1 U | 0.1 UJ | 0.1 | 0.1 | 0.1 | 0.1 | 0.2 | 0.2 |
| Nickel | 3.6 B | 1.6 B | 1.8 | 1.8 | 1.7 | 1.8 | 1.7 | 96 | 40 |
| Potassium | — | — | — | — | — | 2,400 | 2,060 | | 5,000 |
| Selenium | 8.5 J | 4.0 U | 4.0 R | 4.4 | 4.4 | 4.4 R | 4.4 UJ | 8.5 | 5 |
| Silver | 1.9 B | 5.5 B | 0.5 | 0.4 | 0.4 | 0.4 | 0.4 | 10 | 10 |
| Sodium | — | — | — | — | — | 21,500 | 20,700 | | 5,000 |
| Thallium | 3.7 B | 3.6 UJ | 3.6 | 2.6 UJ | 2.6 | 2.6 UJ | 2.6 | 40 | 10 |
| Vanadium | — | — | — | — | — | 0.8 | 1.6 | | 50 |
| Zinc | 0.8 U | 10.7 B | 0.8 UJ | 1.5 | 13.5 | 0.6 UJ | 0.6 UJ | 86 | 20 |
| Inorganics - Metals and Cyanide | | | | | | | | | |
| (Total) | | | | | | | | | |
| Aluminum | — | — | — | — | — | 17,000 | 9,900 | | |
| Antimony | — | 28.7 B | 3.0 | 6.9 | 3.7 | 3.7 | 5.5 | | |
| Arsenic | — | 73.2 | 38.8 | 2.9 | 13.2 | 20.5 | 12.4 | | |
| Barium | — | 1,120 | 852 | 336 | 493 | 568 | 440 | | |
| Beryllium | — | 3.3 B | 2.5 | 0.1 | 0.3 | 1.2 | 1.1 | | |
| Cadmium | — | 2.0 B | 0.2 UJ | 0.2 | 0.2 | 0.2 | 1.0 | | |
| Calcium | — | — | — | — | — | 378,000 | 309,000 | | |
| Chromium | — | 82.3 | 64.2 | 12.3 | 21.4 J | 27.0 | 16.9 | | |
| Cobalt | — | — | — | — | — | 24.1 | 12.3 | | |
| Copper | — | 138 | 108 J | 16.7 | 32.1 | 52.1 | 39.3 | | |
| Cyanide | 4.0 U | 4.0 UJ | 4.0 | 3.0 | 3.0 | 3.0 | 1.0 | 10 | 10 |
| Iron | — | 123,000 | 94,100 | 13,100 | 27,200 J | 45,400 | 25,300 | | |
| Lead | — | 95.4 J | 100 J | 9.6 J | 26.0 J | 46.0 | 23.9 | | |
| Magnesium | — | — | — | — | — | 115,000 | 83,600 | | |
| Manganese | — | — | — | — | — | 2,940 | 988 | | |
| Mercury | — | 0.2 | 0.2 J | 0.1 | 0.1 | 0.1 | 0.1 | | |
| Nickel | — | 114 | 88.2 | 14.1 | 26.0 | 41.2 | 23.4 | | |
| Potassium | — | — | — | — | — | 5,050 | 3,970 | | |
| Selenium | — | 65.7 | 4.0 R | 4.4 | 4.4 R | 4.4 UJ | 4.4 UJ | | |
| Silver | — | 10.6 | 0.5 | 0.4 | 0.4 | 0.4 | 0.4 | | |
| Sodium | — | — | — | — | — | 22,100 | 21,900 | | |
| Thallium | — | 3.6 UJ | 4.9 J | 2.6 UJ | 2.6 UJ | 2.6 | 2.6 | | |
| Vanadium | — | — | — | — | — | 41.5 | 22.2 | | |
| Zinc | — | 379 | 279 J | 61.5 | 87.8 J | 147 J | 72.9 | | |
| Volatile Organic Compounds (VOCs) | | | | | | | | | |
| | BRL | BRL | BRL | BRL | BRL | BRL | BRL | | |
| Semi-Volatile Organic Compounds (SVOCs) | | | | | | | | | |
| | BRL | BRL | BRL | BRL | BRL | BRL | BRL | | |
| Bis (2-ethylhexyl) phthalate | 12.0 U | 10.0 U | 906 | 10.0 U | 10.0 U | 10.0 U | 10.0 U | 49 | 10 |
| Pesticides / PCBs | | | | | | | | | |
| | BRL | BRL | BRL | BRL | BRL | BRL | BRL | | |

- Notes:
- 1) All results expressed in micrograms per liter (µg/L).
 - 2) Standard Inorganic Data Qualifiers have been used.
 - 3) Yellow shading indicate a detection above the Contract Required Quantitation Limit (CRQL)
 - 4) Bold red letters with a thick outline indicates a detection above the Trigger Level.
 - 5) BRL = Below Report Limit; reported data values have a data qualifier of U, J, or UJ
 - 6) — = Constituent not analyzed.
 - 7) U = Not detected at the listed reporting limit.
 - 8) B = An estimated value above the method detection limit (MDL) or the instrument detection limit (IDL) but below the CRQL.
 - 9) UJ = A value less than the CRQL but greater than the MDL.
 - 10) J = The analyte was positively identified; the associated numerical value is the approximate concentration of analyte in the sample.
 - 11) R = The sample results are rejected due to deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte can not be verified.
 - 12) CRQL = Contract Required Quantitation Limit
 - 13) Samples analyzed for Dissolved Inorganics were field filtered using a 0.45 micron, gravity flow filter.
 - 14) Detailed summary tables which list report limits and qualified data values for each compound analyzed for by the laboratory as well as qualified laboratory reports are available upon request.

**Skinner Landfill
West Chester, Ohio
Groundwater Analysis Summary Table for Monitoring Well GW-07R**

| Sampling Event (All Results Expressed in Units of µg/l) | | | | | | | | | |
|---|------------------|-------------|-------------|--------|-------------------|-------------|----------|---------------|-------|
| Compound | Baseline Results | | | | Quarterly Results | | | TRIGGER LEVEL | CRQL |
| | September-02 | December-02 | February-03 | May-03 | August-03 | November-03 | March-04 | | |
| Inorganics - Metals (Dissolved) ¹³ | | | | | | | | | |
| Aluminum | — | — | — | — | — | 25.8 | 25.8 | | 200 |
| Antimony | 5.6 B | 12.4 B | 3.3 | 3.7 | 3.7 | 3.7 | 3.7 | 60 | 60 |
| Arsenic | 10.6 | 7.7 B | 3.6 | 2.9 | 2.9 | 4.5 | 2.9 | 20 | 10 |
| Barium | 100 B | 123 B | 99.9 | 98.8 | 152 J | 131 | 113 | 1,000 | 200 |
| Beryllium | 0.1 U | 0.1 U | 0.1 | 0.1 | 0.1 | 0.1 | 0.1 | 5 | 5 |
| Cadmium | 0.2 U | 0.2 U | 0.2 UJ | 0.2 | 0.2 | 0.2 | 0.2 | 5 | 5 |
| Calcium | — | — | — | — | — | 229,000 | 185,000 | | 5,000 |
| Chromium | 0.7 U | 2.9 B | 2.7 | 2.0 | 2.9 | 0.8 | 1.1 | 11 | 10 |
| Cobalt | — | — | — | — | — | 1.4 | 0.7 | | 50 |
| Copper | 1.0 U | 1.0 U | 1.0 UJ | 1.2 | 1.2 | 1.2 | 1.2 | 25 | 25 |
| Iron | 1,760 | 1,250 | 241 | 146 | 301 J | 3,580 | 32.9 | 7,000 | 100 |
| Lead | 1.3 U | 1.3 UJ | 1.3 R | 1.5 J | 1.5 | 1.5 | 1.5 | 4.2 | 3 |
| Magnesium | — | — | — | — | — | 33,000 | 26,300 | | 5,000 |
| Manganese | — | — | — | — | — | 849 | 914 J | | 15 |
| Mercury | 0.1 U | 0.1 U | 0.1 UJ | 0.1 | 0.1 | 0.1 | 0.1 | 0.2 | 0.2 |
| Nickel | 3.1 B | 1.9 B | 3.4 | 3.2 | 2.4 | 1.6 | 2.2 | 96 | 40 |
| Potassium | — | — | — | — | — | 3,260 | 2,350 | | 5,000 |
| Selenium | 11.3 J | 4.0 U | 4.0 R | 4.4 | 4.4 | 4.4 R | 4.4 UJ | 8.5 | 5 |
| Silver | 0.9 B | 4.7 B | 0.5 | 0.4 | 0.4 | 0.4 | 0.4 | 10 | 10 |
| Sodium | — | — | — | — | — | 42,200 | 25,200 | | 5,000 |
| Thallium | 10.5 | 3.6 UJ | 3.6 | 2.6 J | 2.6 | 2.6 UJ | 2.6 | 40 | 10 |
| Vanadium | — | — | — | — | — | 0.8 | 0.8 | | 50 |
| Zinc | 10.1 B | 11.3 B | 0.8 UJ | 10.2 | 10.9 | 30.7 J | 0.6 UJ | 86 | 20 |
| Inorganics - Metals and Cyanide (Total) | | | | | | | | | |
| Aluminum | — | — | — | — | — | 3,130 J | 7,810 | | |
| Antimony | — | 14.2 B | 3.0 | 3.7 | 3.7 | 3.7 | 6.6 J | | |
| Arsenic | — | 9.8 B | 3.6 | 2.9 | 14.6 | 5.3 | 6.9 | | |
| Barium | — | 454 | 260.0 | 132 | 699 | 204 | 484 | | |
| Beryllium | — | 0.2 B | 0.1 | 0.1 | 0.3 | 0.1 | 0.8 | | |
| Cadmium | — | 0.2 U | 0.2 UJ | 0.2 | 0.2 | 0.2 | 0.9 | | |
| Calcium | — | — | — | — | — | 246,000 | 281,000 | | |
| Chromium | — | 19.3 | 9.1 | 3.7 | 21.5 J | 4.9 | 12.9 | | |
| Cobalt | — | — | — | — | — | 4.3 | 7.0 | | |
| Copper | — | 21.8 B | 8.2 J | 4.2 | 30.4 | 10.0 | 35.5 | | |
| Cyanide | 4.0 U | 4.0 U | 4.0 | 3.0 | 3.0 | 3.0 | 1.5 | 10.0 | 10.0 |
| Iron | — | 24,800 | 10,200 | 2,380 | 29,000 J | 9,890 | 20,200 | | |
| Lead | — | 9.1 J | 1.3 R | 1.5 UJ | 16.8 J | 5.2 | 9.2 | | |
| Magnesium | — | — | — | — | — | 41,600 | 54,000 | | |
| Manganese | — | — | — | — | — | 969 | 1,590 | | |
| Mercury | — | 0.1 U | 0.1 UJ | 0.1 | 0.1 | 0.1 | 0.1 | | |
| Nickel | — | 21.3 B | 10.4 | 4.7 | 25.3 | 10.5 | 17.8 | | |
| Potassium | — | — | — | — | — | 3,780 | 4,510 | | |
| Selenium | — | 4.0 U | 4.0 R | 4.4 | 4.4 R | 4.4 UJ | 4.4 UJ | | |
| Silver | — | 5.2 B | 0.5 | 0.4 | 0.4 | 0.4 | 0.4 | | |
| Sodium | — | — | — | — | — | 41,200 | 31,200 | | |
| Thallium | — | 3.6 UJ | 3.6 UJ | 2.6 UJ | 2.6 UJ | 2.6 | 2.6 | | |
| Vanadium | — | — | — | — | — | 6.5 | 15.3 J | | |
| Zinc | — | 63.1 | 27.6 J | 50.7 | 90.3 J | 22.7 J | 51.2 | | |
| Volatile Organic Compounds (VOCs) | BRL | BRL | BRL | BRL | BRL | BRL | BRL | | |
| Semi-Volatile Organic Compounds (SVOCs) | BRL | BRL | BRL | BRL | BRL | BRL | BRL | | |
| Pesticides / PCBs | BRL | BRL | BRL | BRL | BRL | BRL | BRL | | |

- Notes:
- 1) All results expressed in micrograms per liter (µg/L).
 - 2) Standard Inorganic Data Qualifiers have been used.
 - 3) Yellow shading indicate a detection above the Contract Required Quantitation Limit (CRQL).
 - 4) Bold red letters with a thick outline indicates a detection above the Trigger Level.
 - 5) BRL = Below Report Limit; reported data values have a data qualifier of U, J, or UJ
 - 6) — = Constituent not analyzed.
 - 7) U = Not detected at the listed reporting limit.
 - 8) B = An estimated value above the method detection limit (MDL) or the instrument detection limit (IDL) but below the CRQL.
 - 9) UJ = A value less than the CRQL but greater than the MDL.
 - 10) J = The analyte was positively identified; the associated numerical value is the approximate concentration of analyte in the sample.
 - 11) R = The sample results are rejected due to deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte can not be verified.
 - 12) CRQL = Contract Required Quantitation Limit
 - 13) Samples analyzed for Dissolved Inorganics were field filtered using a 0.45 micron, gravity flow filter.
 - 14) Detailed summary tables which list report limits and qualified data values for each compound analyzed for by the laboratory as well as qualified laboratory reports are available upon request.

**Skinner Landfill
West Chester, Ohio
Groundwater Analysis Summary Table for Monitoring Well GW-58**

| Sampling Event (All Results Expressed in Units of µg/l) | | | | | | | | | |
|---|------------------|-------------|-------------|--------|-------------------|-------------|----------|---------------|-------|
| Compound | Baseline Results | | | | Quarterly Results | | | TRIGGER LEVEL | CRQL |
| | September-02 | December-02 | February-03 | May-03 | August-03 | November-03 | March-04 | | |
| Inorganics - Metals (Dissolved) ¹³ | | | | | | | | | |
| Aluminum | — | — | — | — | — | 25.8 | 25.8 | | 200 |
| Antimony | 5.5 B | 9.7 B | 3.0 | 4.9 | 3.7 | 3.7 | 3.7 | 60 | 60 |
| Arsenic | 8.2 B | 3.6 U | 3.6 | 2.9 | 3.1 | 6.0 | 3.1 | 20 | 10 |
| Barium | 170 B | 50.9 B | 163 | 158 | 162 J | 228 | 156 | 1,000 | 200 |
| Beryllium | 0.1 U | 0.1 U | 0.1 | 0.1 | 0.1 | 0.1 | 0.1 | 5 | 5 |
| Cadmium | 0.2 U | 0.2 U | 0.2 UJ | 0.2 | 0.2 | 0.2 | 0.2 | 5 | 5 |
| Calcium | — | — | — | — | — | 96,400 | 109,000 | | 5,000 |
| Chromium | 0.7 U | 2.9 B | 2.7 | 2.8 | 2.5 | 0.8 | 1.5 | 11 | 10 |
| Cobalt | — | — | — | — | — | 0.4 | 1.3 | | 50 |
| Copper | 1.0 U | 1.0 U | 1.0 UJ | 1.2 | 1.2 | 1.2 | 2.9 | 25 | 25 |
| Iron | 3,440 | 4.9 U | 249 | 1,140 | 488 J | 2,890 | 209 | 7,000 | 100 |
| Lead | 1.3 U | 1.3 U | 1.3 R | 1.5 UJ | 1.5 | 1.5 | 1.5 | 4.2 | 3 |
| Magnesium | — | — | — | — | — | 32,800 | 32,500 | | 5,000 |
| Manganese | — | — | — | — | — | 354 | 549 | | 15 |
| Mercury | 0.1 U | 0.1 U | 0.1 UJ | 0.1 | 0.1 | 0.1 | 0.1 | 0.2 | 0.2 |
| Nickel | 2.6 B | 1.7 B | 2.2 | 2.4 | 1.2 | 1.3 | 2.6 | 96 | 40 |
| Potassium | — | — | — | — | — | 5,210 | 4,550 | | 5,000 |
| Selenium | 4.0 UJ | 4.0 UJ | 4.0 R | 4.4 | 4.4 | 4.4 R | 4.4 UJ | 8.5 | 5 |
| Silver | 0.5 U | 4.1 B | 0.5 | 0.4 | 0.4 | 0.4 | 0.4 | 10 | 10 |
| Sodium | — | — | — | — | — | 34,400 | 32,400 | | 5,000 |
| Thallium | 6.3 B | 3.6 U | 3.6 | 2.6 UJ | 2.6 | 2.6 UJ | 2.6 | 40 | 10 |
| Vanadium | — | — | — | — | — | 0.8 | 1.6 | | 50 |
| Zinc | 1.6 B | 11.0 B | 0.8 UJ | 19.2 | 7.7 | 0.6 UJ | 0.6 | 86 | 20 |
| Inorganics - Metals and Cyanide (Total) | | | | | | | | | |
| Aluminum | — | — | — | — | — | 41,600 | 12,000 | | |
| Antimony | — | 14.3 B | 3.0 | 5.2 | 3.7 | 3.7 | 5.7 | | |
| Arsenic | — | 17.5 | 17.1 | 3.2 | 20.6 | 32.9 | 11.5 J | | |
| Barium | — | 422 | 540 | 367.0 | 391 | 822 | 284 | | |
| Beryllium | — | 1.0 B | 1.3 | 0.3 | 0.7 | 2.9 | 1 | | |
| Cadmium | — | 0.2 U | 0.2 UJ | 0.2 | 0.2 | 1.8 | 1.5 | | |
| Calcium | — | — | — | — | — | 745,000 | 214,000 | | |
| Chromium | — | 51.3 | 63.1 | 14.9 | 42.6 J | 112 | 28.2 | | |
| Cobalt | — | — | — | — | — | 57.2 | 13.4 | | |
| Copper | — | 47.7 | 42.5 J | 27.8 | 43.2 | 138 | 45.7 | | |
| Cyanide | 4.0 U | 4.0 | 4.0 | 3.0 | 3.0 | 3.0 | 0.5 | 10 | 10 |
| Iron | — | 54,500 | 61,900 | 17,000 | 40,800 J | 129,000 | 32,700 | | |
| Lead | — | 19.8 J | 38.5 UJ | 23.0 J | 26.8 J | 92.7 | 19.5 | | |
| Magnesium | — | — | — | — | — | 148,000 | 56,000 | | |
| Manganese | — | — | — | — | — | 4,200 | 1,300 | | |
| Mercury | — | 0.1 U | 0.1 UJ | 0.1 | 0.1 | 0.1 | 0.1 | | |
| Nickel | — | 56.2 | 66.5 | 20.1 | 50.6 | 124 | 32.1 | | |
| Potassium | — | — | — | — | — | 11,800 | 7,640 | | |
| Selenium | — | 4.0 U | 4.0 R | 4.4 | 4.4 R | 4.4 UJ | 4.4 UJ | | |
| Silver | — | 5.9 B | 0.5 | 0.4 | 0.4 | 1.6 | 0.4 | | |
| Sodium | — | — | — | — | — | 36,900 | 33,500 | | |
| Thallium | — | 3.6 UJ | 3.6 UJ | 2.6 UJ | 2.6 UJ | 2.6 | 4.1 J | | |
| Vanadium | — | — | — | — | — | 74.0 | 23.2 | | |
| Zinc | — | 153 | 164 J | 78.7 | 137 J | 367 J | 81.0 J | | |
| Volatile Organic Compounds (VOCs) | BRL | BRL | BRL | BRL | BRL | BRL | BRL | | |
| Semi-Volatile Organic Compounds (SVOCs) | BRL | BRL | BRL | BRL | BRL | BRL | BRL | | |
| Pesticides / PCBs | BRL | BRL | BRL | BRL | BRL | BRL | BRL | | |

Notes:

- 1) All results expressed in micrograms per liter (µg/L).
- 2) Standard Inorganic Data Qualifiers have been used.
- 3) Yellow shading indicate a detection above the Contract Required Quantitation Limit (CRQL).
- 4) Bold red letters with a thick outline indicates a detection above the Trigger Level.
- 5) BRL = Below Report Limit; reported data values have a data qualifier of U, J, or UJ
- 6) — = Constituent not analyzed.
- 7) U = Not detected at the listed reporting limit.
- 8) B = An estimated value above the method detection limit (MDL) or the instrument detection limit (IDL) but below the CRQL.
- 9) UJ = A value less than the CRQL but greater than the MDL.
- 10) J = The analyte was positively identified; the associated numerical value is the approximate concentration of analyte in the sample.
- 11) R = The sample results are rejected due to deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte can not be verified.
- 12) CRQL = Contract Required Quantitation Limit
- 13) Samples analyzed for Dissolved Inorganics were field filtered using a 0.45 micron, gravity flow filter.
- 14) Detailed summary tables which list report limits and qualified data values for each compound analyzed for by the laboratory as well as qualified laboratory reports are available upon request.

Skinner Landfill
West Chester, Ohio
Groundwater Analysis Summary Table for Monitoring Well GW-59

| | | Sampling Event (All Results Expressed in Units of µg/l) | | | | | | | | |
|--|--------------|---|-------------|--------|-----------|-------------------|----------|---------------|-------|--|
| | | Baseline Results | | | | Quarterly Results | | | | |
| Compound | September-02 | December-02 | February-03 | May-03 | August-03 | November-03 | March-04 | TRIGGER LEVEL | CRQL | |
| <u>Inorganics - Metals (Dissolved)</u>¹³ | | | | | | | | | | |
| Aluminum | — | — | — | — | — | 25.8 | 27.2 | | 200 | |
| Antimony | 9.9 B | 13.1 B | 3.0 | 8.0 | 7.2 | 6.5 | 3.7 | 60 | 60 | |
| Arsenic | 13.0 | 5.9 B | 3.6 | 2.9 | 2.9 | 2.9 | 2.9 | 20 | 10 | |
| Barium | 142 B | 65.6 B | 35.7 | 37.2 | 38.1 J | 40.7 | 21.8 | 1,000 | 200 | |
| Beryllium | 0.1 U | 0.1 U | 0.1 | 0.1 | 0.1 | 0.1 | 0.1 | 5 | 5 | |
| Cadmium | 0.2 U | 0.2 U | 0.2 UJ | 0.2 | 0.2 | 0.2 | 0.2 | 5 | 5 | |
| Calcium | — | — | — | — | — | 261,000 | 239,000 | | 5,000 | |
| Chromium | 9.6 B | 4.6 B | 4.9 | 4.4 | 3.3 | 0.8 | 1.8 | 11 | 10 | |
| Cobalt | — | — | — | — | — | 0.4 | 0.4 | | 50 | |
| Copper | 2.2 B | 1.0 U | 1.0 UJ | 3.5 | 3.0 | 4.0 | 2.1 | 25 | 25 | |
| Iron | 4,900 | 4.9 U | 591 | 14.1 | 14.1 | 14.1 | 28.8 | 7,000 | 100 | |
| Lead | 1.5 B | 1.3 UJ | 1.3 R | 1.5 UJ | 1.5 | 1.5 | 1.5 | 4.2 | 3 | |
| Magnesium | — | — | — | — | — | 59,500 | 49,000 | | 5,000 | |
| Manganese | — | — | — | — | — | 27.3 | 4.5 J | | 15 | |
| Mercury | 0.1 U | 0.1 U | 0.1 UJ | 0.1 | 0.1 | 0.1 | 0.1 | 0.2 | 0.2 | |
| Nickel | 12.3 B | 1.8 B | 3.5 | 3.5 | 3 | 2.3 | 2.6 | 96 | 40 | |
| Potassium | — | — | — | — | — | 29,800 | 32,800 | | 5,000 | |
| Selenium | 15.8 J | 18.2 | 4.0 R | 4.4 | 4.4 | 4.4 R | 4.4 UJ | 8.5 | 5 | |
| Silver | 2.7 B | 7.2 B | 0.5 | 0.4 | 0.4 | 0.4 | 0.4 | 10 | 10 | |
| Sodium | — | — | — | — | — | 186,000 | 166,000 | | 5,000 | |
| Thallium | 8.8 B | 3.6 UJ | 3.6 UJ | 2.6 UJ | 2.6 | 2.6 UJ | 3.1 J | 40 | 10 | |
| Vanadium | — | — | — | — | — | 0.8 | 1.4 | | 50 | |
| Zinc | 7.9 B | 7.2 B | 0.8 J | 1.2 | 18.4 | 0.6 UJ | 3.1 J | 86 | 20 | |
| <u>Inorganics - Metals and Cyanide</u> | | | | | | | | | | |
| <u>(Total)</u> | | | | | | | | | | |
| Aluminum | — | — | — | — | — | 3,710 | 816 | | | |
| Antimony | — | 19.5 B | 3.0 | 3.7 | 3.7 | 3.7 | 4.7 | | | |
| Arsenic | — | 36.5 | 3.6 | 2.9 | 3.6 | 4.3 | 2.9 | | | |
| Barium | — | 215 | 202 | 55.2 | 62.1 | 213 | 55.0 | | | |
| Beryllium | — | 1.9 B | 0.1 | 0.1 | 0.1 | 0.1 | 0.2 | | | |
| Cadmium | — | 0.7 B | 0.2 UJ | 0.2 | 0.2 | 0.2 | 0.2 | | | |
| Calcium | — | — | — | — | — | 281,000 | 243,000 | | | |
| Chromium | — | 82.4 | 18.8 | 5.4 | 8.3 J | 19.1 | 5.5 | | | |
| Cobalt | — | — | — | — | — | 7.4 | 2.1 | | | |
| Copper | — | 45.0 | 1.0 UJ | 3.6 | 6.0 | 11.9 | 10.1 | | | |
| Cyanide | 4.0 U | 4.0 | 4.0 | 3.0 | 3.0 | 3.0 | 1.0 | 10 | 10 | |
| Iron | — | 79,700 | 9,810 | 1,390 | 2,240 J | 12,900 | 3,020 | | | |
| Lead | — | 36.7 J | 1.3 R | 1.5 UJ | 5.7 J | 10.0 | 1.5 | | | |
| Magnesium | — | — | — | — | — | 62,400 | 51,500 | | | |
| Manganese | — | — | — | — | — | 923 | 224 | | | |
| Mercury | — | 0.1 B | 0.1 UJ | 0.1 | 0.1 | 0.1 | 0.1 | | | |
| Nickel | — | 77.2 | 16.0 | 4.8 | 6.5 | 20.0 | 6.7 | | | |
| Potassium | — | — | — | — | — | 31,900 | 32,500 | | | |
| Selenium | — | 21.6 | 4.0 R | 4.4 | 4.4 R | 4.4 UJ | 4.4 UJ | | | |
| Silver | — | 7.6 B | 0.5 | 0.4 | 0.4 | 0.4 | 0.4 | | | |
| Sodium | — | — | — | — | — | 180,000 | 162,000 | | | |
| Thallium | — | 3.6 UJ | 3.6 UJ | 2.6 UJ | 2.6 UJ | 2.6 | 2.6 | | | |
| Vanadium | — | — | — | — | — | 5.9 | 2.3 | | | |
| Zinc | — | 238 | 18.9 J | 16.3 | 21.2 J | 36.3 J | 7.9 J | | | |
| <u>Volatile Organic Compounds (VOCs)</u> | | | | | | | | | | |
| 1,1-Dichloroethane | BRL | BRL | BRL | BRL | BRL | BRL | BRL | | 10 | |
| <u>Semi-Volatile Organic Compounds (SVOCs)</u> | | | | | | | | | | |
| | BRL | BRL | BRL | BRL | BRL | BRL | BRL | | | |
| <u>Pesticides / PCBs</u> | | | | | | | | | | |
| | BRL | BRL | BRL | BRL | BRL | BRL | BRL | | | |

Notes:

- 1) All results expressed in micrograms per liter (µg/L).
- 2) Standard Inorganic Data Qualifiers have been used.
- 3) Yellow shading indicate a detection above the Contract Required Quantitation Limit (CRQL).
- 4) Bold red letters with a thick outline indicates a detection above the Trigger Level.
- 5) BRL = Below Report Limit; reported data values have a data qualifier of U, J, or UJ
- 6) — = Constituent not analyzed.
- 7) U = Not detected at the listed reporting limit.
- 8) B = An estimated value above the method detection limit (MDL) or the instrument detection limit (IDL) but below the CRQL.
- 9) UJ = A value less than the CRQL but greater than the MDL.
- 10) J = The analyte was positively identified; the associated numerical value is the approximate concentration of analyte in the sample.
- 11) R = The sample results are rejected due to deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte can not be verified.
- 12) CRQL = Contract Required Quantitation Limit
- 13) Samples analyzed for Dissolved Inorganics were field filtered using a 0.45 micron, gravity flow filter.
- 14) Detailed summary tables which list report limits and qualified data values for each compound analyzed for by the laboratory as well as qualified laboratory reports are available upon request.

Skinner Landfill
West Chester, Ohio
Groundwater Analysis Summary Table for Monitoring Well GW-60

| Sampling Event (All Results Expressed in Units of µg/l) | | | | | | | | | |
|---|---------------------|-------------|---------------------|--------|---------------------|-------------|---------------------|---------------|-------|
| Compound | Baseline Results | | | | Quarterly Results | | | TRIGGER LEVEL | CRQL |
| | September-02 | December-02 | February-03 | May-03 | August-03 | November-03 | March-04 | | |
| Inorganics - Metals (Dissolved)¹³ | Insufficient Volume | | Insufficient Volume | | Insufficient Volume | | Insufficient Volume | | |
| Aluminum | — | — | — | — | — | 25.8 | 25.8 | | 200 |
| Antimony | — | 11.3 B | — | 6.7 | — | 3.7 | 5.1 | 60 | 60 |
| Arsenic | — | 3.6 U | — | 2.9 | — | 2.9 | 2.9 | 20 | 10 |
| Barium | — | 43 B | — | 60.9 | — | 28.7 | 27.1 | 1,000 | 200 |
| Beryllium | — | 0.1 U | — | 0.1 | — | 0.1 | 0.2 | 5 | 5 |
| Cadmium | — | 0.2 U | — | 0.2 | — | 0.2 | 0.2 | 5 | 5 |
| Calcium | — | — | — | — | — | 100,000 | 309,000 | | 5,000 |
| Chromium | — | 4.6 B | — | 2.7 | — | 0.8 | 2.6 | 11 | 10 |
| Cobalt | — | — | — | — | — | 0.4 | 0.4 | | 50 |
| Copper | — | 1.0 U | — | 3.0 | — | 4.2 | 4.0 | 25 | 25 |
| Iron | — | 1.3 UJ | — | 14.1 | — | 14.1 | 14.1 | 7,000 | 100 |
| Lead | — | 0.1 U | — | 1.5 UJ | — | 1.5 | 1.5 | 4.2 | 3 |
| Magnesium | — | — | — | — | — | 20,100 | 88,200 | | 5,000 |
| Manganese | — | — | — | — | — | 2.4 | 0.5 J | | 15 |
| Mercury | — | 3.4 B | — | 0.1 | — | 0.1 | 0.1 | 0.2 | 0.2 |
| Nickel | — | 13.3 | — | 1.1 | — | 0.7 | 2.4 | 96 | 40 |
| Potassium | — | — | — | — | — | 6,970 | 6,480 | | 5,000 |
| Selenium | — | 5.8 B | — | 4.4 | — | 4.4 R | 4.4 UJ | 8.5 | 5 |
| Silver | — | 3.6 UJ | — | 0.4 | — | 0.4 | 0.4 | 10 | 10 |
| Sodium | — | — | — | — | — | 201,000 | 46,000 | | 5,000 |
| Thallium | — | 10.4 B | — | 2.6 UJ | — | 2.6 UJ | 2.6 | 40 | 10 |
| Vanadium | — | — | — | — | — | 0.8 | 0.8 | | 50 |
| Zinc | — | 10.4 B | — | 4.5 | — | 0.6 UJ | 0.6 UJ | 86 | 20 |
| Inorganics - Metals and Cyanide (Total) | | | | | | | | | |
| Aluminum | — | — | — | — | — | 13,400 J | 32,500 | | |
| Antimony | — | 8.4 B | — | 3.7 | — | 3.7 | 9.7 | | |
| Arsenic | — | 5.7 B | — | 2.9 | — | 11.7 | 17.0 J | | |
| Barium | — | 88.5 B | — | 73.1 | — | 89.8 | 129 | | |
| Beryllium | — | 0.1 U | — | 0.1 | — | 0.9 | 2.5 | | |
| Cadmium | — | 0.2 U | — | 0.2 | — | 0.2 | 2.8 | | |
| Calcium | — | — | — | — | — | 158,000 | 492,000 | | |
| Chromium | — | 7.3 B | — | 10.9 | — | 33.2 | 59.6 | | |
| Cobalt | — | — | — | — | — | 16.6 | 36.1 | | |
| Copper | — | 1.0 U | — | 7.5 | — | 29.3 | 54.5 | | |
| Cyanide | — | 4.0 U | — | 3.0 | — | 3.0 | — | 10 | 10 |
| Iron | — | 2,780 | — | 7830 | — | 31,300 | 74,200 | | |
| Lead | — | 1.3 UJ | — | 1.5 UJ | — | 28.2 | 40.4 | | |
| Magnesium | — | — | — | — | — | 32,500 | 112,000 | | |
| Manganese | — | — | — | — | — | 555 | 1,410 | | |
| Mercury | — | 0.1 U | — | 0.1 | — | 0.1 | 0.1 | | |
| Nickel | — | 7.4 B | — | 7.8 | — | 31.6 | 67.3 | | |
| Potassium | — | — | — | — | — | 9,290 | 11,800 | | |
| Selenium | — | 14.3 | — | 4.4 | — | 4.4 UJ | 4.4 UJ | | |
| Silver | — | 5.0 B | — | 0.4 | — | 0.4 | 0.4 | | |
| Sodium | — | — | — | — | — | 212,000 | 44,600 | | |
| Thallium | — | 3.6 UJ | — | 2.6 UJ | — | 2.6 | 11 J | | |
| Vanadium | — | — | — | — | — | 23.2 | 51.2 | | |
| Zinc | — | 28.3 | — | 34.2 | — | 135 J | 180 J | | |
| Volatile Organic Compounds (VOCs) | BRL | BRL | BRL | BRL | BRL | BRL | BRL | | |
| Semi-Volatile Organic Compounds (SVOCs) | — | BRL | BRL | BRL | — | BRL | BRL | | |
| Pesticides / PCBs | — | BRL | — | BRL | — | BRL | BRL | | |

Notes:

- 1) All results expressed in micrograms per liter (µg/L).
- 2) Standard Inorganic Data Qualifiers have been used.
- 3) Yellow shading indicate a detection above the Contract Required Quantitation Limit (CRQL).
- 4) Bold red letters with a thick outline indicates a detection above the Trigger Level.
- 5) BRL = Below Report Limit; reported data values have a data qualifier of U, J, or UJ
- 6) — = Constituent not analyzed.
- 7) U = Not detected at the listed reporting limit.
- 8) B = An estimated value above the method detection limit (MDL) or the instrument detection limit (IDL) but below the CRQL.
- 9) UJ = A value less than the CRQL but greater than the MDL.
- 10) — = The analyte was positively identified; the associated numerical value is the approximate concentration of analyte in the sample.
- 11) — = The sample results are rejected due to deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte can not be verified.
- 12) CRQL = Contract Required Quantitation Limit
- 13) Samples analyzed for Dissolved Inorganics were field filtered using a 0.45 micron, gravity flow filter.
- 14) Detailed summary tables which list report limits and qualified data values for each compound analyzed for by the laboratory as well as qualified laboratory reports are available upon request.

**Skinner Landfill
West Chester, Ohio
Groundwater Analysis Summary Table for Monitoring Well GW-61**

| Compound | Sampling Event (All Results Expressed in Units of µg/l) | | | | | | | TRIGGER LEVEL | CRQL |
|--|---|-------------|-------------|--------|-------------------|-------------|----------|------------------|-------|
| | Baseline Results | | | | Quarterly Results | | | | |
| | September-02 | December-02 | February-03 | May-03 | August-03 | November-03 | March-04 | | |
| <u>Inorganics - Metals (Dissolved)¹³</u> | | | | | | | | | |
| Aluminum | — | — | — | — | — | 25.8 | 25.8 | | 200 |
| Antimony | 6.3 B | 6.7 B | 3.0 | 7.3 | 3.7 | 3.7 | 4.5 | 60 | 60 |
| Arsenic | 10.6 | 13.7 | 3.6 | 2.9 | 4.7 | 7.5 | 2.9 | 20 | 10 |
| Barium | 104 B | 98.2 B | 64.7 | 67.7 | 77.7 J | 83.3 | 39.4 | 1,000 | 200 |
| Beryllium | 0.1 U | 0.1 U | 0.1 | 0.1 | 0.1 | 0.1 | 0.1 | 5 | 5 |
| Cadmium | 0.2 U | 0.2 U | 0.2 UJ | 0.2 | 0.2 | 0.2 | 0.3 | 5 | 5 |
| Calcium | — | — | — | — | — | 191,000 | 191,000 | | 5,000 |
| Chromium | 0.7 U | 4.4 B | 1.6 | 3.3 | 3.8 | 0.8 | 1.1 | 11 | 10 |
| Cobalt | — | — | — | — | — | 2.0 | 1.4 | | 50 |
| Copper | 1.0 U | 1.0 U | 1.0 UJ | 1.2 | 1.2 | 1.2 | 8.0 | 25 | 25 |
| Iron | 2,770 | 12,500 | 3,270 | 1,940 | 6,100 J | 5,100 | 187 | 7,000 | 100 |
| Lead | 1.3 U | 1.3 UJ | 1.3 R | 1.5 UJ | 1.5 | 1.5 | 1.5 | 4.2 | 3 |
| Magnesium | — | — | — | — | — | 35,700 | 29,100 | | 5,000 |
| Manganese | — | — | — | — | — | 866 | 485 | | 15 |
| Mercury | 0.1 U | 0.1 U | 0.1 UJ | 0.1 | 0.1 | 0.1 | 0.1 | 0.2 | 0.2 |
| Nickel | 3.1 B | 5.0 B | 2.4 | 4.2 | 4.0 | 4.0 | 4.2 | 96 | 40 |
| Potassium | — | — | — | — | — | 10,100 | 6,990 | | 5,000 |
| Selenium | 5.3 J | 4.0 U | 4.0 R | 4.4 | 4.4 | 4.4 R | 4.4 UJ | 8.5 | 5 |
| Silver | 0.7 B | 3.4 B | 0.5 | 0.4 | 0.4 | 0.4 | 0.4 | 10 | 10 |
| Sodium | — | — | — | — | — | 28,300 | 27,900 | | 5,000 |
| Thallium | 5.6 B | 3.6 UJ | 3.6 | 2.6 UJ | 2.6 | 2.6 UJ | 2.6 | 40 | 10 |
| Vanadium | — | — | — | — | — | 0.8 | 1.2 | | 50 |
| Zinc | 4.1 B | 18.5 B | 0.8 UJ | 7.4 | 13.1 | 4.8 J | 0.6 | 86 | 20 |
| <u>Inorganics - Metals and Cyanide (Total)</u> | | | | | | | | | |
| Aluminum | — | — | — | — | — | 1,080 | 452 | | |
| Antimony | — | 6.7 B | 3.0 | 4.5 | 3.7 | 3.7 | 4.8 | | |
| Arsenic | — | 13.7 | 3.6 | 3.2 | 17.9 | 3.7 | 2.9 | | |
| Barium | — | 98.2 B | 84.4 | 69.5 | 202 | 91.3 | 44.1 | | |
| Beryllium | — | 0.1 U | 0.1 | 0.1 | 0.2 | 0.1 | 0.2 | | |
| Cadmium | — | 0.2 U | 0.2 UJ | 0.2 | 0.2 | 0.2 | 0.3 | | |
| Calcium | — | — | — | — | — | 190,000 | 187,000 | | |
| Chromium | — | 4.4 B | 5.6 | 2.9 | 23.2 J | 2.1 | 1.9 | | |
| Cobalt | — | — | — | — | — | 3.3 | 1.7 | | |
| Copper | — | 1.0 U | 1.0 UJ | 1.2 | 28.2 | 4.2 | 22.2 | | |
| Cyanide | 4.0 U | 4.0 U | 4.0 | 3.0 | 3.0 | 3.0 | 0.5 | 10 | 10 |
| Iron | — | 12,500 | 8,720 | 2,330 | 33,400 J | 8,640 | 2,430 | | |
| Lead | — | 1.3 UJ | 1.3 R | 1.5 UJ | 19.7 J | 1.6 | 22.1 | | |
| Magnesium | — | — | — | — | — | 37,500 | 30,000 | | |
| Manganese | — | — | — | — | — | 922 | 527 | | |
| Mercury | — | 0.1 U | 0.1 UJ | 0.1 | 0.1 | 0.1 | 0.1 | | |
| Nickel | — | 5.0 B | 6.6 | 4.1 | 29.5 | 7.6 | 4.3 | | |
| Potassium | — | — | — | — | — | 9,430 | 6,950 | | |
| Selenium | — | 4.0 U | 4.0 R | 4.4 | 4.4 R | 4.4 UJ | 4.4 UJ | | |
| Silver | — | 3.4 B | 0.5 | 0.4 | 0.4 | 0.4 | 0.4 | | |
| Sodium | — | — | — | — | — | 27,700 | 27,000 | | |
| Thallium | — | 3.6 UJ | 3.8 UJ | 2.6 UJ | 2.6 UJ | 2.6 | 2.6 | | |
| Vanadium | — | — | — | — | — | 0.8 | 2.1 | | |
| Zinc | — | 18.5 B | 9.9 J | 18.3 | 96.7 J | 13.8 J | 7.3 J | | |
| <u>Volatile Organic Compounds (VOCs)</u> | | | | | | | | | |
| Carbon disulfide | 1.0 U | 1.0 U | 1.0 U | 1.0 U | 1.4 | 1.0 U | 1.0 U | | 10 |
| <u>Semi-Volatile Organic Compounds (SVOCs)</u> | | | | | | | | | |
| Bis (2-ethylhexyl) phthalate | 12.0 U | 57.0 J | 10.0 U | 10.0 U | 40.6 | 10.0 U | 10.0 U | 49 | 10 |
| <u>Pesticides / PCBs</u> | | | | | | | | | |
| | BRL | BRL | BRL | BRL | BRL | BRL | BRL | | |

Notes:

- 1) All results expressed in micrograms per liter (µg/L).
- 2) Standard Inorganic Data Qualifiers have been used.
- 3) Yellow shading indicate a detection above the Contract Required Quantitation Limit (CRQL).
- 4) Bold red letters with a thick outline indicates a detection above the Trigger Level.
- 5) BRL = Below Report Limit; reported data values have a data qualifier of U, J, or UJ
- 6) — = Constituent not analyzed.
- 7) U = Not detected at the listed reporting limit.
- 8) B = An estimated value above the method detection limit (MDL) or the instrument detection limit (IDL) but below the CRQL.
- 9) UJ = A value less than the CRQL but greater than the MDL.
- 10) J = The analyte was positively identified; the associated numerical value is the approximate concentration of analyte in the sample.
- 11) R = The sample results are rejected due to deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte can not be verified.
- 12) CRQL = Contract Required Quantitation Limit
- 13) Samples analyzed for Dissolved Inorganics were field filtered using a 0.45 micron, gravity flow filter.
- 14) Detailed summary tables which list report limits and qualified data values for each compound analyzed for by the laboratory as well as qualified laboratory reports are available upon request.

**Skinner Landfill
West Chester, Ohio
Groundwater Analysis Summary Table for Monitoring Well GW-62A**

| | Sampling Event (All Results Expressed in Units of µg/l) | | | | | | | | |
|---|---|-------------|-------------|--------|-------------------|-------------|----------|---------------|-------|
| | Baseline Results | | | | Quarterly Results | | | | |
| Compound | September-02 | December-02 | February-03 | May-03 | August-03 | November-03 | March-04 | TRIGGER LEVEL | CRQL |
| Inorganics - Metals (Dissolved) ¹³ | | | | | | | | | |
| Aluminum | — | — | — | — | — | 25.8 | 25.8 | | 200 |
| Antimony | 7.3 B | 12.7 B | 3.0 | 9.6 | 3.7 | 3.7 | 3.7 | 60 | 60 |
| Arsenic | 4.9 B | 3.6 U | 3.6 | 2.9 | 2.9 | 2.9 | 2.9 | 20 | 10 |
| Barium | 174 B | 157 B | 162 | 146 | 145 J | 126 | 111 | 1,000 | 200 |
| Beryllium | 0.1 U | 0.1 U | 0.1 | 0.1 | 0.1 | 0.1 | 0.1 | 5 | 5 |
| Cadmium | 0.2 U | 0.2 U | 0.2 UJ | 0.2 | 0.2 | 0.2 | 0.2 | 5 | 5 |
| Calcium | — | — | — | — | — | 123,000 | 122,000 | | 5,000 |
| Chromium | 1.5 B | 4.1 B | 3.5 | 3.5 | 3.5 | 0.8 | 2.1 | 11 | 10 |
| Cobalt | — | — | — | — | — | 0.4 | 0.5 | | 50 |
| Copper | 1.1 B | 5.2 B | 1.0 UJ | 4.4 | 2.4 | 2.7 | 1.2 | 25 | 25 |
| Iron | 6.2 B | 4.9 U | 317 | 14.1 | 14.1 | 14.1 | 14.1 | 7,000 | 100 |
| Lead | 1.3 U | 1.3 UJ | 1.3 R | 1.5 UJ | 1.5 | 1.5 | 1.5 | 4.2 | 3 |
| Magneisum | — | — | — | — | — | 49,200 | 48,700 | | 5,000 |
| Manganese | — | — | — | — | — | 51.4 | 164 J | | 15 |
| Mercury | 0.1 UJ | 0.1 U | 0.1 UJ | 0.1 | 0.1 | 0.1 | 0.1 | 0.2 | 0.2 |
| Nickel | 3.2 B | 2.7 B | 1.8 | 2.1 | 0.7 | 0.9 | 1.8 | 96 | 40 |
| Potassium | — | — | — | — | — | 10,800 | 11,100 | | 5,000 |
| Selenium | 4.0 U | 4.0 U | 4.0 R | 4.4 | 4.4 | 4.4 R | 4.4 UJ | 8.5 | 5 |
| Silver | 0.5 U | 6.0 B | 0.5 | 0.5 | 0.4 | 0.4 | 0.4 | 10 | 10 |
| Sodium | — | — | — | — | — | 120,000 | 120,000 | | 5,000 |
| Thallium | 7.9 B | 3.6 UJ | 3.6 | 2.6 UJ | 2.6 | 2.6 UJ | 2.6 | 40 | 10 |
| Vanadium | — | — | — | — | — | 0.8 | 1.6 | | 50 |
| Zinc | 0.8 U | 11.4 B | 0.8 UJ | 7.4 | 11.0 | 0.9 J | 0.6 UJ | 86 | 20 |
| Inorganics - Metals and Cyanide | | | | | | | | | |
| (Total) | | | | | | | | | |
| Aluminum | — | — | — | — | — | 24,100 | 13,200 | | |
| Antimony | — | 18.4 B | 3.0 | 4.7 | 3.7 | 3.7 | 6.1 | | |
| Arsenic | — | 17.0 | 20.7 | 8.7 | 18.3 | 17.7 | 8.3 J | | |
| Barium | — | 471 | 1,170 | 615 | 800 | 633 | 361 | | |
| Beryllium | — | 0.5 B | 1.7 | 0.8 | 1.1 | 1.5 | 1.1 | | |
| Cadmium | — | 0.2 U | 0.2 UJ | 0.2 | 0.2 | 1.1 | 1.6 | | |
| Calcium | — | — | — | — | — | 618,000 | 337,000 | | |
| Chromium | — | 38.6 | 85.5 | 49.3 | 68.6 J | 49.5 | 29.6 | | |
| Cobalt | — | — | — | — | — | 33.5 | 15.6 | | |
| Copper | — | 42.3 | 76.1 UJ | 45.2 | 68 | 72.8 | 42.7 | | |
| Cyanide | 4.0 U | 4.0 U | 4.0 | — | 3.0 | 3.0 | 1.0 | 10.0 | 10.0 |
| Iron | — | 34,000 | 85,100 | 51,500 | 65,400 J | 60,800 | 35,000 | | |
| Lead | — | 33.3 J | 68.0 UJ | 33.6 J | 65.2 J | 72.8 | 39.5 | | |
| Magneisum | — | — | — | — | — | 137,000 | 88,000 | | |
| Manganese | — | — | — | — | — | 3,380 | 1,460 | | |
| Mercury | — | 0.1 U | 0.1 UJ | 0.1 | 0.1 | 0.1 | 0.1 | | |
| Nickel | — | 38.1 B | 86.4 | 53.4 | 75.7 | 64.3 | 35.4 | | |
| Potassium | — | — | — | — | — | 15,100 | 13,900 | | |
| Selenium | — | 13.7 | 4.0 R | 4.4 | 4.4 R | 4.4 UJ | 4.4 UJ | | |
| Silver | — | 5.4 B | 0.5 | 0.4 | 0.4 | 0.4 | 0.4 | | |
| Sodium | — | — | — | — | — | 121,000 | 123,000 | | |
| Thallium | — | 3.6 UJ | 3.6 UJ | 2.6 UJ | 2.6 UJ | 2.6 | 6.9 J | | |
| Vanadium | — | — | — | — | — | 40.5 | 23.0 | | |
| Zinc | — | 119 | 242 J | 179 | 234 J | 181 J | 101 J | | |
| Volatile Organic Compounds (VOCs) | BRL | BRL | BRL | BRL | BRL | BRL | BRL | | |
| Semi-Volatile Organic Compounds (SVOCs) | BRL | BRL | BRL | BRL | BRL | BRL | BRL | | |
| Pesticides / PCBs | BRL | BRL | BRL | BRL | BRL | BRL | BRL | | |

Notes:

- 1) All results expressed in micrograms per liter (µg/L).
- 2) Standard Inorganic Data Qualifiers have been used.
- 3) Yellow shading indicate a detection above the Contract Required Quantitation Limit (CRQL).
- 4) Bold red letters with a thick outline indicates a detection above the Trigger Level.
- 5) BRL = Below Report Limit; reported data values have a data qualifier of U, J, or UJ
- 6) — = Constituent not analyzed.
- 7) U = Not detected at the listed reporting limit.
- 8) B = An estimated value above the method detection limit (MDL) or the instrument detection limit (IDL) but below the CRQL.
- 9) UJ = A value less than the CRQL but greater than the MDL.
- 10) J = The analyte was positively identified; the associated numerical value is the approximate concentration of analyte in the sample.
- 11) R = The sample results are rejected due to deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte can not be verified.
- 12) CRQL = Contract Required Quantitation Limit
- 13) Samples analyzed for Dissolved Inorganics were field filtered using a 0.45 micron, gravity flow filter.
- 14) Detailed summary tables which list report limits and qualified data values for each compound analyzed for by the laboratory as well as qualified laboratory reports are available upon request.

Skinner Landfill
West Chester, Ohio
Groundwater Analysis Summary Table for Monitoring Well GW-62B

| Compound | Sampling Event (All Results Expressed in Units of µg/l) | | | | | | | TRIGGER LEVEL | CRQL |
|---|---|-------------|-------------|---------------------|-------------------|-------------|-------------|------------------|------|
| | Baseline Results | | | | Quarterly Results | | | | |
| | September-02 | December-02 | February-03 | May-03 | August-03 | November-03 | March-04 | | |
| <u>Inorganics - Metals (Dissolved)</u> ¹³ | Well is Dry | Well is Dry | Well is Dry | Insufficient Volume | Well is Dry | Well is Dry | Well is Dry | | |
| <u>Inorganics - Metals and Cyanide (Total)</u> | — | — | — | — | — | — | — | | |
| <u>Volatile Organic Compounds (VOCs)</u> | — | — | — | BRL | — | — | — | | |
| Benzene | — | — | — | 3.00 | — | — | — | 5 | 10 |
| <u>Semi-Volatile Organic Compounds (SVOCs)</u> | — | — | — | BRL | — | — | — | | |
| Bis (2-Chloroethyl) ether | — | — | — | 24.5 | — | — | — | 13.6 | 10 |
| <u>Pesticides / PCBs</u> | — | — | — | BRL | — | — | — | | |

Notes:

- 1) All results expressed in micrograms per liter (µg/L).
- 2) Standard Inorganic Data Qualifiers have been used.
- 3) Yellow shading indicate a detection above the Contract Required Quantitation Limit (CRQL)
- 4) Bold red letters with a thick outline indicates a detection above the Trigger Level.
- 5) BRL = Below Report Limit; reported data values have a data qualifier of U, J, or UJ
- 6) — = No Sample Available (Well Dry)
- 7) U = Not detected at the listed reporting limit.
- 8) B = An estimated value above the method detection limit (MDL) or the instrument detection limit (IDL) but below the CRQL.
- 9) UJ = A value less than the CRQL but greater than the MDL.
- 10) J = The analyte was positively identified; the associated numerical value is the approximate concentration of analyte in the sample.
- 11) R = The sample results are rejected due to deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte can not be verified.
- 12) CRQL = Contract Required Quantitation Limit
- 13) Samples analyzed for Dissolved Inorganics were field filtered using a 0.45 micron, gravity flow filter.
- 14) Detailed summary tables which list report limits and qualified data values for each compound analyzed for by the laboratory as well as qualified laboratory reports are available upon request.

**Skinner Landfill
West Chester, Ohio
Groundwater Analysis Summary Table for Monitoring Well GW-63**

| | | Sampling Event (All Results Expressed in Units of µg/l) | | | | | | | |
|---|--------------|---|-------------|--------|-----------|-------------------|----------|---------------|-------|
| | | Baseline Results | | | | Quarterly Results | | | |
| Compound | September-02 | December-02 | February-03 | May-03 | August-03 | November-03 | March-04 | TRIGGER LEVEL | CRQL |
| Inorganics - Metals (Dissolved) ¹³ | | | | | | | | | |
| Aluminum | — | — | — | — | — | 25.8 | 25.8 | | 200 |
| Antimony | 11.4 B | 13.8 B | 3.0 | 8.9 | 7.1 | 3.7 | 3.8 | 60 | 60 |
| Arsenic | 15.9 | 9.5 B | 3.6 | 2.9 | 2.9 | 5.4 | 2.9 | 20 | 10 |
| Barium | 97.5 B | 76.2 B | 72.2 | 50.1 | 58.8 J | 68.6 | 20.1 | 1,000 | 200 |
| Beryllium | 0.1 U | 0.1 U | 0.1 | 0.1 | 0.1 | 0.1 | 0.2 | 5 | 5 |
| Cadmium | 0.2 U | 0.2 U | 0.2 UJ | 0.2 | 0.2 | 0.2 | 0.2 | 5 | 5 |
| Calcium | — | — | — | — | — | 278,000 | 295,000 | | 5,000 |
| Chromium | 1.2 B | 3 B | 3.6 | 3.4 | 3.2 | 0.8 | 1.8 | 11 | 10 |
| Cobalt | — | — | — | — | — | 4.1 | 1.1 | | 50 |
| Copper | 1.0 U | 1.0 U | 1.0 UJ | 1.2 | 1.2 | 1.2 | 2.0 J | 25 | 25 |
| Iron | 592 | 758 | 622 | 297 | 1,890 J | 1,150 | 21.4 | 7,000 | 100 |
| Lead | 1.3 U | 1.3 UJ | 1.3 R | 1.5 UJ | 1.5 | 1.5 | 1.5 | 4.2 | 3 |
| Magnesium | — | — | — | — | — | 61,000 | 67,000 | | 5,000 |
| Manganese | — | — | — | — | — | 2,600 | 271 J | | 15 |
| Mercury | 0.1 UJ | 0.1 U | 0.1 UJ | 0.1 | 0.1 | 0.1 | 0.1 | 0.2 | 0.2 |
| Nickel | 12.8 B | 7.4 B | 6.6 | 5.9 | 8.3 | 6.9 | 3.2 | 96 | 40 |
| Potassium | — | — | — | — | — | 11,600 | 5,210 | | 5,000 |
| Selenium | 15.3 | 8.7 | 4.0 R | 4.4 | 4.4 | 4.4 R | 4.4 UJ | 8.5 | 5 |
| Silver | 2.0 B | 3.6 B | 0.5 | 0.4 | 0.4 | 0.4 | 0.4 | 10 | 10 |
| Sodium | — | — | — | — | — | 72,100 | 46,100 | | 5,000 |
| Thallium | 15.9 | 3.6 UJ | 3.6 | 2.6 UJ | 2.6 | 2.6 UJ | 4.6 J | 40 | 10 |
| Vanadium | — | — | — | — | — | 0.8 | 0.8 | | 50 |
| Zinc | 0.8 U | 7.8 B | 0.8 UJ | 12.9 | 10.8 | 3.7 J | 0.6 UJ | 86 | 20 |
| Inorganics - Metals and Cyanide (Total) | | | | | | | | | |
| Aluminum | — | — | — | — | — | 10,500 | 26,600 | | |
| Antimony | — | 20.2 B | 3.0 | 4.6 | 3.7 | 3.7 | 5.7 | | |
| Arsenic | — | 22.8 | 30.4 | 12.6 | 12.6 | 9.3 | 17.1 J | | |
| Barium | — | 234 | 390 | 178 | 217 | 147 | 186 | | |
| Beryllium | — | 1.0 B | 2.3 | 0.9 | 0.9 | 0.6 | 2.1 | | |
| Cadmium | — | 0.2 U | 0.2 UJ | 0.2 | 0.2 | 0.2 | 2.5 | | |
| Calcium | — | — | — | — | — | 465,000 | 465,000 | | |
| Chromium | — | 39.2 | 70.8 | 30.8 | 36.1 J | 13.7 | 38.2 | | |
| Cobalt | — | — | — | — | — | 17.5 | 28.3 | | |
| Copper | — | 35.7 | 77.8 J | 29.3 | 33.0 | 17.4 | 69.2 | | |
| Cyanide | 4.0 U | 4.0 U | 4.0 | 3.0 | 3.0 | 3.0 | 0.5 | 10 | 10 |
| Iron | — | 55,500 | 109,000 | 44,100 | 53,900 J | 25,800 | 63,200 | | |
| Lead | — | 21.2 J | 70.2 UJ | 28.5 J | 36 J | 23.4 | 41.0 | | |
| Magnesium | — | — | — | — | — | 96,100 | 111,000 | | |
| Manganese | — | — | — | — | — | 4,090 | 2,570 | | |
| Mercury | — | 0.1 U | 0.1 UJ | 0.1 | 0.1 | 0.1 | 0.1 | | |
| Nickel | — | 56.6 | 104 | 47.5 | 59.6 | 31.0 | 58.1 | | |
| Potassium | — | — | — | — | — | 31,500 | 9,320 | | |
| Selenium | — | 27.8 | 4.0 R | 4.4 | 4.4 R | 4.4 UJ | 4.4 UJ | | |
| Silver | — | 7.3 B | 0.5 | 0.4 | 0.4 | 0.4 | 0.4 | | |
| Sodium | — | — | — | — | — | 73,600 | 45,000 | | |
| Thallium | — | 3.6 U | 3.6 UJ | 2.6 UJ | 2.6 UJ | 2.6 | 8.5 J | | |
| Vanadium | — | — | — | — | — | 17.8 | 43 | | |
| Zinc | — | 160 | 267 J | 129 | 182 J | 66.3 J | 176 J | | |
| Volatile Organic Compounds (VOCs) | | | | | | | | | |
| Acetone | 5.0 U | 5.0 R | 5.0 U | 78.0 J | 5.0 U | 5.0 U | 5.0 R | | 10 |
| Carbon disulfide | 1.0 U | 1.0 U | 1.0 U | 1.0 U | 1.3 | 1.0 U | 1.0 U | | 10 |
| Semi-Volatile Organic Compounds (SVOCs) | | | | | | | | | |
| butylbenzylphthalate | BRL | BRL | BRL | BRL | BRL | BRL | BRL | | |
| | | | | | | | 0.61 J | 10 | 10 |
| Pesticides / PCBs | | | | | | | | | |
| | BRL | BRL | BRL | BRL | BRL | BRL | BRL | | |

Notes:

- 1) All results expressed in micrograms per liter (µg/L).
- 2) Standard Inorganic Data Qualifiers have been used.
- 3) Yellow shading indicate a detection above the Contract Required Quantitation Limit (CRQL)
- 4) Bold red letters with a thick outline indicates a detection above the Trigger Level.
- 5) BRL = Below Report Limit; reported data values have a data qualifier of U, J, or UJ
- 6) — = Constituent not analyzed.
- 7) U = Not detected at the listed reporting limit.
- 8) B = An estimated value above the method detection limit (MDL) or the instrument detection limit (IDL) but below the CRQL.
- 9) UJ = A value less than the CRQL but greater than the MDL.
- 10) J = The analyte was positively identified; the associated numerical value is the approximate concentration of analyte in the sample.
- 11) R = The sample results are rejected due to deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte can not be verified.
- 12) CRQL = Contract Required Quantitation Limit
- 13) Samples analyzed for Dissolved Inorganics were field filtered using a 0.45 micron, gravity flow filter.
- 14) Detailed summary tables which list report limits and qualified data values for each compound analyzed for by the laboratory as well as qualified laboratory reports are available upon request.

**Skinner Landfill
West Chester, Ohio
Groundwater Analysis Summary Table for Monitoring Well GW-64**

| Sampling Event (All Results Expressed in Units of µg/l) | | | | | | | | | |
|---|------------------|-------------|-------------|--------|-------------------|-------------|----------|---------------|-------|
| | Baseline Results | | | | Quarterly Results | | | | |
| Compound | September-02 | December-02 | February-03 | May-03 | August-03 | November-03 | March-04 | TRIGGER LEVEL | CRQL |
| Inorganics - Metals (Dissolved) ¹³ | | | | | | | | | |
| Aluminum | — | — | — | — | — | 25.8 | 25.8 | | 200 |
| Antimony | 5.9 B | 12.3 B | 3.0 | 5.3 | 4.2 | 3.7 | 3.7 | 60 | 60 |
| Arsenic | 9.2 B | 7.3 B | 3.6 | 2.9 | 2.9 | 2.9 | 2.9 | 20 | 10 |
| Barium | 34.6 B | 53.5 B | 31.0 | 28.9 | 31.5 J | 44.6 | 28.3 | 1,000 | 200 |
| Beryllium | 0.1 U | 0.1 U | 0.1 | 0.1 | 0.1 | 0.1 | 0.1 | 5 | 5 |
| Cadmium | 0.2 U | 0.2 U | 0.2 UJ | 0.2 | 0.2 | 0.2 | 0.2 | 5 | 5 |
| Calcium | — | — | — | — | — | 185,000 | 176,000 | | 5,000 |
| Chromium | 1.2 B | 20.4 | 4.9 | 6.5 | 3.8 | 0.8 | 1.7 | 11 | 10 |
| Cobalt | — | — | — | — | — | 0.5 | 1.8 | | 50 |
| Copper | 1.8 B | 10.4 B | 1.0 UJ | 4.9 | 4.5 | 3.4 | 1.2 | 25 | 25 |
| Iron | 4.9 U | 19,500 | 63.2 R | 52.6 | 14.1 | 14.1 | 14.1 | 7,000 | 100 |
| Lead | 1.3 U | 3.7 J | 1.3 UJ | 1.5 UJ | 1.5 | 1.5 | 1.5 | 4.2 | 3 |
| Magnesium | — | — | — | — | — | 61,800 | 56,700 | | 5,000 |
| Manganese | — | — | — | — | — | 292 | 1,170 J | | 15 |
| Mercury | 0.1 U | 0.1 U | 0.1 | 0.1 | 0.1 | 0.1 | 0.1 | 0.2 | 0.2 |
| Nickel | 10.9 B | 33.6 B | 12.8 R | 13.4 | 9.4 | 5.2 | 7.8 | 96 | 40 |
| Potassium | — | — | — | — | — | 12,300 | 12,900 | | 5,000 |
| Selenium | 6.6 J | 4.6 B | 4.0 | 4.4 | 4.4 | 4.4 R | 4.4 UJ | 8.5 | 5 |
| Silver | 0.8 B | 5.7 B | 0.5 | 0.4 | 0.4 | 0.4 | 0.4 | 10 | 10 |
| Sodium | — | — | — | — | — | 67,600 | 53,900 | | 5,000 |
| Thallium | 3.9 B | 3.6 UJ | 3.6 | 2.6 UJ | 2.6 | 2.6 UJ | 2.6 | 40 | 10 |
| Vanadium | — | — | — | — | — | 0.8 | 0.8 | | 50 |
| Zinc | 0.8 U | 64.5 | 0.8 UJ | 9.1 | 16.1 | 2.6 J | 0.6 UJ | 86 | 20 |
| Inorganics - Metals and Cyanide | | | | | | | | | |
| (Total) | | | | | | | | | |
| Aluminum | — | — | — | — | — | 18,700 J | 3,080 | | |
| Antimony | — | 12.3 B | 3.0 | 5.6 | 3.7 | 3.7 | 4.9 | | |
| Arsenic | — | 7.3 B | 4.3 | 2.9 | 6.2 | 10.8 | 2.9 | | |
| Barium | — | 53.5 B | 59.5 | 47.9 | 58.3 | 95.9 | 37.1 | | |
| Beryllium | — | 0.1 U | 0.5 | 0.1 | 0.2 | 1.0 | 0.3 | | |
| Cadmium | — | 0.2 U | 0.2 UJ | 0.2 | 0.2 | 0.2 | 0.2 | | |
| Calcium | — | — | — | — | — | 311,000 | 213,000 | | |
| Chromium | — | 20.4 | 22.9 | 14.8 | 22.4 J | 29.4 | 7.0 | | |
| Cobalt | — | — | — | — | — | 23.1 | 5.4 | | |
| Copper | — | 10.4 B | 2.2 J | 9.6 | 16.0 | 16.3 | 11.3 | | |
| Cyanide | 4.0 U | 4.0 U | 4.0 | 3.0 | 3.0 | 3.0 | 1.3 | 10 | 10 |
| Iron | — | 19,500 | 25,900 | 14,700 | 24,300 J | 42,900 | 7,520 | | |
| Lead | — | 3.7 J | 8.1 UJ | 1.7 J | 9.7 J | 20.0 | 1.5 | | |
| Magnesium | — | — | — | — | — | 77,300 | 66,000 | | |
| Manganese | — | — | — | — | — | 2,390 | 1,650 | | |
| Mercury | — | 0.1 U | 0.1 UJ | 0.1 | 0.1 | 0.1 | 0.1 | | |
| Nickel | — | 33.6 B | 38.3 | 26.1 | 33.9 | 46.0 | 16.4 | | |
| Potassium | — | — | — | — | — | 14,700 | 15,000 | | |
| Selenium | — | 4.6 B | 4.0 R | 4.4 | 4.4 R | 4.4 UJ | 4.4 UJ | | |
| Silver | — | 5.7 B | 0.5 | 0.4 | 0.4 | 0.4 | 0.4 | | |
| Sodium | — | — | — | — | — | 68,300 | 59,800 | | |
| Thallium | — | 3.6 UJ | 3.6 UJ | 2.6 UJ | 2.6 UJ | 2.6 | 2.6 | | |
| Vanadium | — | — | — | — | — | 27.3 | 5.3 | | |
| Zinc | — | 64.5 | 51.3 J | 69.1 | 73.6 J | 114 J | 13.6 J | | |
| Volatile Organic Compounds (VOCs) | BRL | BRL | BRL | BRL | BRL | BRL | BRL | | |
| Semi-Volatile Organic Compounds (SVOCs) | BRL | BRL | BRL | BRL | BRL | BRL | BRL | | |
| Bis (2-ethylhexyl) phthalate | 12.0 U | 10.0 U | 10.0 U | 10.0 U | 10.0 | 10.0 U | 10.0 U | 49 | 10 |
| Pesticides / PCBs | BRL | BRL | BRL | BRL | BRL | BRL | BRL | | |

- Notes:
- 1) All results expressed in micrograms per liter (µg/L).
 - 2) Standard Inorganic Data Qualifiers have been used.
 - 3) Yellow shading indicate a detection above the Contract Required Quantitation Limit (CRQL)
 - 4) Bold red letters with a thick outline indicates a detection above the Trigger Level.
 - 5) BRL = Below Report Limit; reported data values have a data qualifier of U, J, or UJ
 - 6) — = Constituent not analyzed.
 - 7) U = Not detected at the listed reporting limit.
 - 8) B = An estimated value above the method detection limit (MDL) or the instrument detection limit (IDL) but below the CRQL.
 - 9) UJ = A value less than the CRQL but greater than the MDL.
 - 10) J = The analyte was positively identified; the associated numerical value is the approximate concentration of analyte in the sample.
 - 11) R = The sample results are rejected due to deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte can not be verified.
 - 12) CRQL = Contract Required Quantitation Limit
 - 13) Samples analyzed for Dissolved Inorganics were field filtered using a 0.45 micron, gravity flow filter.
 - 14) Detailed summary tables which list report limits and qualified data values for each compound analyzed for by the laboratory as well as qualified laboratory reports are available upon request.

Skinner Landfill
West Chester, Ohio
Groundwater Analysis Summary Table for Monitoring Well GW-65

| Sampling Event (All Results Expressed in Units of µg/l) | | | | | | | | | |
|---|---------------------|---------------------|---------------------|---------------------|---------------------|---------------------|---------------------|---------------|------|
| Compound | Baseline Results | | | | Quarterly Results | | | TRIGGER LEVEL | CRQL |
| | September-02 | December-02 | February-03 | May-03 | August-03 | November-03 | March-04 | | |
| Inorganics - Metals (Dissolved)¹³ | Insufficient Volume | Insufficient Volume | Insufficient Volume | Insufficient Volume | Insufficient Volume | Insufficient Volume | Insufficient Volume | | |
| Antimony | — | — | — | — | — | — | — | 60 | 60 |
| Arsenic | — | — | — | — | — | — | — | 10 | 10 |
| Barium | — | — | — | — | — | — | — | 1,000 | 200 |
| Beryllium | — | — | — | — | — | — | — | 5 | 5 |
| Cadmium | — | — | — | — | — | — | — | 5 | 5 |
| Chromium | — | — | — | — | — | — | — | 11 | 10 |
| Copper | — | — | — | — | — | — | — | 25 | 25 |
| Iron | — | — | — | — | — | — | — | 5,000 | 100 |
| Lead | — | — | — | — | — | — | — | 4.2 | 3 |
| Mercury | — | — | — | — | — | — | — | 0.2 | 0.2 |
| Nickel | — | — | — | — | — | — | — | 96 | 40 |
| Selenium | — | — | — | — | — | — | — | 5 | 5 |
| Silver | — | — | — | — | — | — | — | 10 | 10 |
| Thallium | — | — | — | — | — | — | — | 40 | 10 |
| Zinc | — | — | — | — | — | — | — | 86 | 20 |
| Inorganics - Metals and Cyanide (Total) | | | | | | | | | |
| Antimony | — | — | — | — | — | — | — | | |
| Arsenic | — | — | — | — | — | — | — | | |
| Barium | — | — | — | — | — | — | — | | |
| Beryllium | — | — | — | — | — | — | — | | |
| Cadmium | — | — | — | — | — | — | — | | |
| Chromium | — | — | — | — | — | — | — | | |
| Copper | — | — | — | — | — | — | — | | |
| Cyanide | — | — | — | — | — | — | — | 10 | 10 |
| Iron | — | — | — | — | — | — | — | | |
| Lead | — | — | — | — | — | — | — | | |
| Mercury | — | — | — | — | — | — | — | | |
| Nickel | — | — | — | — | — | — | — | | |
| Selenium | — | — | — | — | — | — | — | | |
| Silver | — | — | — | — | — | — | — | | |
| Thallium | — | — | — | — | — | — | — | | |
| Zinc | — | — | — | — | — | — | — | | |
| Volatile Organic Compounds (VOCs) | BRL | BRL | BRL | BRL | BRL | BRL | BRL | | |
| Semi-Volatile Organic Compounds (SVOCs) | BRL | BRL | BRL | BRL | BRL | BRL | — | | |
| 4-Nitrophenol | 50.0 U | 29.8 UJ | 10.0 U | 10.0 U | 10.0 U | 10.0 U | — | 150 | 25 |
| Bis (2-ethylhexyl) phthalate | 20.0 U | 11.9 U | 10.0 U | 10.0 U | 10.0 U | 10.0 U | — | 49 | 10 |
| Pesticides / PCBs | — | — | — | — | — | — | — | | |

Notes:

- 1) All results expressed in micrograms per liter (µg/L).
- 2) Standard Inorganic Data Qualifiers have been used.
- 3) Yellow shading indicate a detection above the Contract Required Quantitation Limit (CRQL)
- 4) Bold red letters with a thick outline indicates a detection above the Trigger Level.
- 5) BRL = Below Report Limit; reported data values have a data qualifier of U, J, or UJ
- 6) — = No Sample Available (Well Dry)
- 7) U = Not detected at the listed reporting limit.
- 8) B = An estimated value above the method detection limit (MDL) or the instrument detection limit (IDL) but below the CRQL.
- 9) UJ = A value less than the CRQL but greater than the MDL.
- 10) J = The analyte was positively identified; the associated numerical value is the approximate concentration of analyte in the sample.
- 11) R = The sample results are rejected due to deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte can not be verified.
- 12) CRQL = Contract Required Quantitation Limit
- 13) Samples analyzed for Dissolved Inorganics were field filtered using a 0.45 micron, gravity flow filter.
- 14) Detailed summary tables which list report limits and qualified data values for each compound analyzed for by the laboratory as well as qualified laboratory reports are available upon request.

**Skinner Landfill
West Chester, Ohio**
Groundwater Analysis Summary Table for Creek Surface Water Sample Location SW-50

| | | Sampling Event (All Results Expressed in Units of µg/l) | | | | | | | | |
|---|--------------|---|-------------|--------|-----------|-------------------|-----------|---------------|-------|--|
| | | Baseline Results | | | | Quarterly Results | | | | |
| Compound | September-02 | December-02 | February-03 | May-03 | August-03 | November-03 | March-04 | TRIGGER LEVEL | CRQL | |
| Inorganics - Metals (Dissolved) ¹³ | | | | | | | | | | |
| Aluminum | — | — | — | — | — | 25.8 | 25.8 | | 200 | |
| Antimony | 3.9 B | 9.7 B | 3.0 U | 9.7 B | 3.7 | 3.7 | 3.7 | 60 | 60 | |
| Arsenic | 3.6 U | 3.6 U | 3.6 U | 3.6 U | 7.1 | 2.9 | 2.9 | 20 | 10 | |
| Barium | 69.3 B | 50.9 B | 57.7 B | 50.9 B | 55.2 | 40.0 | 35.8 | 1,000 | 200 | |
| Beryllium | 0.1 U | 0.1 U | 0.1 U | 0.1 U | 0.1 | 0.1 | 0.1 | 5 | 5 | |
| Cadmium | 0.2 U | 0.2 U | 0.2 UJ | 0.2 U | 0.2 | 0.2 | 0.2 | 5 | 5 | |
| Calcium | — | — | — | — | — | 84,100 | 103,000 | | 5,000 | |
| Chromium | 0.7 U | 2.9 B | 2.1 B | 2.9 B | 1.8 | 0.8 | 1.6 | 11 | 10 | |
| Cobalt | — | — | — | — | — | 0.5 | 0.4 | | 50 | |
| Copper | 1.0 U | 1.0 U | 1.0 UJ | 1.0 U | 3.1 | 4.4 | 4.3 | 25 | 25 | |
| Iron | 56.0 B | 4.9 U | 129 | 4.9 U | 14.1 | 14.1 | 14.1 | 7,000 | 100 | |
| Lead | 1.3 U | 1.3 U | 1.3 R | 1.3 U | 1.5 | 1.5 | 1.5 | 4.2 | 3 | |
| Magnesium | — | — | — | — | — | 23,400 | 29,700 | | 5,000 | |
| Manganese | — | — | — | — | — | 3.8 | 30.0 J | | 15 | |
| Mercury | 0.1 U | 0.1 U | 0.1 U | 0.1 U | 0.1 | 0.1 | 0.1 | 0.2 | 0.2 | |
| Nickel | 0.8 B | 1.7 B | 0.8 B | 1.7 B | 0.7 | 0.7 | 1.0 | 96 | 40 | |
| Potassium | — | — | — | — | — | 3,840 | 2,980 | | 5,000 | |
| Selenium | 4.0 UJ | 4.0 UJ | 4.0 R | 4.0 UJ | 4.4 | 4.4 R | 4.4 R | 8.5 | 5 | |
| Silver | 0.7 B | 4.1 B | 0.5 U | 4.1 B | 0.4 | 0.4 | 0.4 | 10 | 10 | |
| Sodium | — | — | — | — | — | 32,100 | 59,200 | | 5,000 | |
| Thallium | 4.0 B | 3.6 U | 4.3 J | 3.6 U | 6.8 J | 2.6 | 2.6 | 40 | 10 | |
| Vanadium | — | — | — | — | — | 0.8 | 2.2 | | 50 | |
| Zinc | 1.3 B | 11.0 B | 0.8 UJ | 11.0 B | 13.8 | 0.6 UJ | 0.6 | 86 | 20 | |
| Inorganics - Metals and Cyanide (Total) | | | | | | | | | | |
| Aluminum | — | — | — | — | — | 25.8 | 62,300 J | | | |
| Antimony | — | 9.3 B | 3.0 U | 9.3 B | 3.7 | 3.7 | 7.3 J | | | |
| Arsenic | — | 3.6 U | 3.6 U | 3.6 U | 2.9 | 3.4 | 50.7 J | | | |
| Barium | — | 50.7 B | 58.1 B | 50.7 B | 55.4 | 41.3 | 499 J | | | |
| Beryllium | — | 0.1 U | 0.1 U | 0.1 U | 0.1 | 0.1 | 4.9 | | | |
| Cadmium | — | 0.2 U | 0.2 UJ | 0.2 U | 0.2 | 0.2 | 5.0 | | | |
| Calcium | — | — | — | — | — | 86,400 | 427,000 J | | | |
| Chromium | — | 3.2 B | 2.3 B | 3.2 B | 1.7 | 0.8 | 72.6 J | | | |
| Cobalt | — | — | — | — | — | 0.4 | 59.7 | | | |
| Copper | — | 4.4 B | 5.7 J | 4.4 B | 2.5 | 4.4 | 131 J | | | |
| Cyanide | 4.0 U | 4.0 U | 4.0 U | 4.0 U | 3.0 | 3.0 | 0.8 | 10 | 10 | |
| Iron | — | 4.9 U | 141 | 4.9 U | 23.5 | 69.2 | 124,000 J | | | |
| Lead | — | 1.3 U | 1.3 R | 1.3 U | 1.5 | 1.5 | 122 J | | | |
| Magnesium | — | — | — | — | — | 23,900 | 80,300 | | | |
| Manganese | — | — | — | — | — | 5.8 | 5,690 | | | |
| Mercury | — | 0.1 U | 0.1 U | 0.1 U | 0.1 | 0.1 | 0.1 | | | |
| Nickel | — | 1.3 B | 0.5 U | 1.3 B | 1.5 | 0.7 | 116 J | | | |
| Potassium | — | — | — | — | — | 3,990 | 12,200 | | | |
| Selenium | — | 4.0 U | 4.0 R | 4.0 U | 4.4 UJ | 4.4 R | 4.4 R | | | |
| Silver | — | 4.3 B | 0.5 U | 4.3 B | 0.4 | 0.4 | 0.4 | | | |
| Sodium | — | — | — | — | — | 33,000 | 60,200 | | | |
| Thallium | — | 3.6 UJ | 3.6 UJ | 3.6 UJ | 5.5 J | 2.6 | 2.6 | | | |
| Vanadium | — | — | — | — | — | 0.8 | 105 | | | |
| Zinc | — | 13.1 B | 2.0 J | 13.1 B | 10.9 | 1.3 J | 490 J | | | |
| Volatile Organic Compounds (VOCs) | BRL | BRL | BRL | BRL | BRL | BRL | BRL | | | |
| Semi-Volatile Organic Compounds (SVOCs) | BRL | BRL | BRL | BRL | BRL | BRL | BRL | | | |
| Fluoranthene | | | | | | | 0.84 J | 10 | 10 | |
| Phenanthrene | | | | | | | 0.79 J | 10 | 10 | |
| Pyrene | | | | | | | 0.67 J | 10 | 10 | |
| Diethylphthalate | | | | | | 2.17 J | 10.0 U | 10 | 10 | |
| Pesticides / PCBs | BRL | BRL | BRL | BRL | BRL | BRL | BRL | | | |

Notes:

- 1) All results expressed in micrograms per liter (µg/L).
- 2) Standard Inorganic Data Qualifiers have been used.
- 3) Yellow shading indicate a detection above the Contract Required Quantitation Limit (CRQL).
- 4) Bold red letters with a thick outline indicates a detection above the Trigger Level.
- 5) BRL = Below Report Limit; reported data values have a data qualifier of U, J, or UJ
- 6) — = Constituent not analyzed.
- 7) U = Not detected at the listed reporting limit.
- 8) B = An estimated value above the method detection limit (MDL) or the instrument detection limit (IDL) but below the CRQL.
- 9) UJ = A value less than the CRQL but greater than the MDL.
- 10) J = The analyte was positively identified; the associated numerical value is the approximate concentration of analyte in the sample.
- 11) R = The sample results are rejected due to deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte can not be verified.
- 12) CRQL = Contract Required Quantitation Limit
- 13) Samples analyzed for Dissolved Inorganics were field filtered using a .45 micron, gravity flow filter.
- 14) Detailed summary tables which list report limits and qualified data values for each compound analyzed for by the laboratory as well as qualified laboratory reports are available upon request.

Skinner Landfill
West Chester, Ohio
Groundwater Analysis Summary Table for Creek Surface Water Sample Location SW-51

| Sampling Event (All Results Expressed in Units of µg/l) | | | | | | | | | |
|---|------------------|-------------|-------------|--------|-------------------|-------------|-----------|---------------|-------|
| Compound | Baseline Results | | | | Quarterly Results | | | TRIGGER LEVEL | CRQL |
| | September-02 | December-02 | February-03 | May-03 | August-03 | November-03 | March-04 | | |
| Inorganics - Metals (Dissolved) ¹³ | | | | | | | | | |
| Aluminum | — | — | — | — | — | 25.8 | 25.8 | | 200 |
| Antimony | 5.3 B | 8.7 B | 3.0 U | 8.7 B | 3.7 | 3.7 | 3.7 | 60 | 60 |
| Arsenic | 5.8 U | 3.6 U | 3.6 U | 3.6 U | 5.3 | 3.2 | 2.9 | 20 | 10 |
| Barium | 60.2 B | 54.8 B | 59.4 B | 54.8 B | 49.9 | 42.2 | 36.7 | 1,000 | 200 |
| Beryllium | 0.1 U | 0.1 U | 0.1 U | 0.1 U | 0.1 | 0.1 | 0.1 | 5 | 5 |
| Cadmium | 0.2 U | 0.2 U | 0.2 UJ | 0.2 U | 0.2 | 0.2 | 0.2 | 5 | 5 |
| Calcium | — | — | — | — | — | 88,800 | 106,000 | | 5,000 |
| Chromium | 0.7 U | 2.7 B | 2.5 B | 2.7 B | 0.8 | 0.8 | 1.7 | 11 | 10 |
| Cobalt | — | — | — | — | — | 0.4 | 0.4 | | 50 |
| Copper | 1.0 U | 1.0 U | 1.0 UJ | 1.0 U | 2.2 | 5.1 | 3.0 | 25 | 25 |
| Iron | 6.4 B | 4.9 U | 11.9 B | 4.9 U | 14.1 | 14.1 | 14.1 | 7,000 | 100 |
| Lead | 1.3 U | 1.3 U | 1.3 R | 1.3 U | 1.5 | 1.5 | 1.5 | 4.2 | 3 |
| Magnesium | — | — | — | — | — | 24,700 | 31,500 | | 5,000 |
| Manganese | — | — | — | — | — | 4.6 | 29.9 J | | 15 |
| Mercury | 0.1 U | 0.1 U | 0.1 U | 0.1 U | 0.1 | 0.1 | 0.1 | 0.2 | 0.2 |
| Nickel | 0.7 B | 1.7 B | 0.5 U | 1.7 B | 0.7 | 0.7 | 2.4 | 96 | 40 |
| Potassium | — | — | — | — | — | 3,910 | 2,160 | | 5,000 |
| Selenium | 4.0 UJ | 4.0 UJ | 4.0 R | 4.0 UJ | 4.4 | 4.4 R | 4.4 R | 8.5 | 5 |
| Silver | 0.5 U | 3.3 B | 0.5 U | 3.3 B | 0.4 | 0.4 | 0.4 | 10 | 10 |
| Sodium | — | — | — | — | — | 34,000 | 60,200 | | 5,000 |
| Thallium | 3.6 U | 3.6 U | 3.6 UJ | 3.6 U | 10.6 | 2.6 | 2.6 | 40 | 10 |
| Vanadium | — | — | — | — | — | 0.8 | 2.4 | | 50 |
| Zinc | 0.8 U | 26.8 | 0.8 UJ | 26.8 | 13.9 | 0.6 UJ | 3.4 | 86 | 20 |
| Inorganics - Metals and Cyanide (Total) | | | | | | | | | |
| Aluminum | — | — | — | — | — | 25.8 | 9,250 J | | |
| Antimony | — | 7.8 B | 3.0 U | 7.8 B | 3.7 | 3.7 | 3.7 UJ | | |
| Arsenic | — | 3.6 U | 3.6 U | 3.6 U | 2.9 | 2.9 | 11.1 J | | |
| Barium | — | 51.8 B | 59.3 B | 51.8 B | 50.9 | 42.6 | 112 J | | |
| Beryllium | — | 0.1 U | 0.1 U | 0.1 U | 0.1 | 0.1 | 0.7 | | |
| Cadmium | — | 0.2 U | 0.2 UJ | 0.2 U | 0.2 | 0.2 | 0.5 | | |
| Calcium | — | — | — | — | — | 86,700 | 153,000 J | | |
| Chromium | — | 2.7 B | 2.8 B | 2.7 B | 1.7 | 0.8 | 12.7 J | | |
| Cobalt | — | — | — | — | — | 0.4 | 7.8 | | |
| Copper | — | 1.0 U | 1.0 UJ | 1.0 U | 1.9 | 3.2 | 22.0 J | | |
| Cyanide | 4.0 U | 4.0 U | 4.0 U | 4.0 U | 3.0 | 3.0 | 0.6 | 10 | 10 |
| Iron | — | 10.0 B | 74.8 B | 10.0 B | 49.5 | 83.8 | 17,800 J | | |
| Lead | — | 1.3 U | 1.3 R | 1.3 U | 1.5 | 1.5 | 17.3 J | | |
| Magnesium | — | — | — | — | — | 23,900 | 38,900 J | | |
| Manganese | — | — | — | — | — | 6.5 | 685 J | | |
| Mercury | — | 0.1 U | 0.1 U | 0.1 U | 0.1 | 0.1 | 0.1 | | |
| Nickel | — | 1.1 B | 0.5 U | 1.1 B | 0.7 | 0.7 | 16.4 J | | |
| Potassium | — | — | — | — | — | 3,820 | 4,470 | | |
| Selenium | — | 4.0 U | 4.0 R | 4.0 U | 4.4 UJ | 4.4 R | 4.4 R | | |
| Silver | — | 3.0 B | 0.5 U | 3.0 B | 0.4 | 0.4 | 0.4 | | |
| Sodium | — | — | — | — | — | 32,800 | 61,800 | | |
| Thallium | — | 3.6 UJ | 3.6 U | 3.6 UJ | 5.9 J | 2.6 | 2.6 | | |
| Vanadium | — | — | — | — | — | 0.8 | 18.7 J | | |
| Zinc | — | 12.7 B | 7.8 J | 12.7 B | 9.7 | 0.6 UJ | 52.9 J | | |
| Volatile Organic Compounds (VOCs) | BRL | BRL | BRL | BRL | BRL | BRL | BRL | | |
| Semi-Volatile Organic Compounds (SVOCs) | BRL | BRL | BRL | BRL | BRL | BRL | BRL | | |
| Pesticides / PCBs | BRL | BRL | BRL | BRL | BRL | BRL | BRL | | |

Notes:

- 1) All results expressed in micrograms per liter (µg/L).
- 2) Standard Inorganic Data Qualifiers have been used.
- 3) Yellow shading indicate a detection above the Contract Required Quantitation Limit (CRQL).
- 4) Bold red letters with a thick outline indicates a detection above the Trigger Level.
- 5) BRL = Below Report Limit; reported data values have a data qualifier of U, J, or UJ
- 6) — = Constituent not analyzed.
- 7) U = Not detected at the listed reporting limit.
- 8) B = An estimated value above the method detection limit (MDL) or the instrument detection limit (IDL) but below the CRQL.
- 9) UJ = A value less than the CRQL but greater than the MDL.
- 10) J = The analyte was positively identified; the associated numerical value is the approximate concentration of analyte in the sample.
- 11) R = The sample results are rejected due to deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte can not be verified.
- 12) CRQL = Contract Required Quantitation Limit
- 13) Samples analyzed for Dissolved Inorganics were field filtered using a .45 micron, gravity flow filter.
- 14) Detailed summary tables which list report limits and qualified data values for each compound analyzed for by the laboratory as well as qualified laboratory reports are available upon request.

**Skinner Landfill
West Chester, Ohio
Groundwater Analysis Summary Table for Creek Surface Water Sample Location SW-52**

| Sampling Event (All Results Expressed in Units of µg/l) | | | | | | | | | |
|---|--------------|-------------|-------------|--------|-------------------|-------------|-----------|------------------|-------|
| Baseline Results | | | | | Quarterly Results | | | TRIGGER LEVEL | CRQL |
| Compound | September-02 | December-02 | February-03 | May-03 | August-03 | November-03 | March-04 | | |
| Inorganics - Metals (Dissolved)¹³ | | | | | | | | | |
| Aluminum | — | — | — | — | — | 25.8 | 25.8 | | 200 |
| Antimony | 4.6 B | 6.5 B | 3.0 U | 6.5 B | 3.7 | 3.7 | 3.7 | 60 | 60 |
| Arsenic | 5.0 B | 3.6 U | 3.6 U | 3.6 U | 4.9 | 2.9 | 2.9 | 20 | 10 |
| Barium | 64.3 B | 52.8 B | 57.7 B | 52.8 B | 54.2 | 40.7 | 40.7 | 1,000 | 200 |
| Beryllium | 0.1 U | 0.1 U | 0.1 U | 0.1 U | 0.1 | 0.1 | 0.1 | 5 | 5 |
| Cadmium | 0.2 U | 0.2 U | 0.2 UJ | 0.2 U | 0.2 | 0.2 | 0.2 | 5 | 5 |
| Calcium | — | — | — | — | — | 86,000 | 107,000 | | 5,000 |
| Chromium | 0.8 B | 2.8 B | 3.5 B | 2.8 B | 2.1 | 0.8 | 1.3 | 11 | 10 |
| Cobalt | — | — | — | — | — | 0.4 | 0.4 | | 50 |
| Copper | 1.0 U | 1.0 U | 1.0 UJ | 1.0 U | 2.3 | 3.7 | 2.0 | 25 | 25 |
| Iron | 4.9 U | 4.9 U | 575 | 4.9 U | 14.1 | 14.1 | 14.1 | 7,000 | 100 |
| Lead | 1.3 U | 1.3 U | 1.3 R | 1.3 U | 1.5 | 1.5 | 1.5 | 4.2 | 3 |
| Magnesium | — | — | — | — | — | 23,500 | 30,700 | | 5,000 |
| Manganese | — | — | — | — | — | 5.0 | 1.5 J | | 15 |
| Mercury | 0.1 U | 0.1 U | 0.1 U | 0.1 U | 0.1 | 0.1 | 0.1 | 0.2 | 0.2 |
| Nickel | 0.5 U | 1.2 B | 1.5 B | 1.2 B | 1.1 | 0.7 | 0.7 | 96 | 40 |
| Potassium | — | — | — | — | — | 3,720 | 1,900 | | 5,000 |
| Selenium | 4.0 UJ | 4.0 UJ | 4.0 R | 4.0 UJ | 4.4 | 4.4 R | 4.4 R | 8.5 | 5 |
| Silver | 0.5 U | 3.5 B | 0.5 U | 3.5 B | 0.4 | 0.4 | 0.4 | 10 | 10 |
| Sodium | — | — | — | — | — | 32,900 | 61,400 | | 5,000 |
| Thallium | 3.6 U | 3.6 U | 3.6 UJ | 3.6 U | 6.2 J | 2.6 | 2.6 | 40 | 10 |
| Vanadium | — | — | — | — | — | 0.8 | 1.9 | | 50 |
| Zinc | 0.8 U | 15.8 B | 0.8 UJ | 15.8 B | 26.5 | 1.5 J | 4.5 | 86 | 20 |
| <u>Inorganics - Metals and Cyanide (Total)</u> | | | | | | | | | |
| Aluminum | — | — | — | — | — | 25.8 | 26.8 J | | |
| Antimony | — | 5.9 B | 3.0 U | 5.9 B | 3.7 | 3.7 | 3.7 UJ | | |
| Arsenic | — | 3.6 U | 3.6 U | 3.6 U | 2.9 | 2.9 | 2.9 | | |
| Barium | — | 53.2 B | 58.4 B | 53.2 B | 54.4 | 41.2 | 41.7 J | | |
| Beryllium | — | 0.1 U | 0.1 U | 0.1 U | 0.1 | 0.1 | 0.1 | | |
| Cadmium | — | 0.2 U | 0.2 UJ | 0.2 U | 0.2 | 0.2 | 0.2 | | |
| Calcium | — | — | — | — | — | 84,800 | 108,000 J | | |
| Chromium | — | 3.7 B | 2.4 B | 3.7 B | 1.5 | 0.8 | 1.7 J | | |
| Cobalt | — | — | — | — | — | 0.4 | 0.5 | | |
| Copper | — | 1.0 U | 1.0 UJ | 1.0 U | 1.7 | 3.3 | 3.2 J | | |
| Cyanide | 4.0 U | 4.0 U | 4.0 B | 4.0 U | 3.0 | 3.0 | 0.5 J | 10 | 10 |
| Iron | — | 28.2 B | 76.7 R | 28.2 B | 45.9 | 79.6 | 67.8 J | | |
| Lead | — | 1.3 U | 1.3 R | 1.3 U | 1.5 | 1.5 | 1.5 | | |
| Magnesium | — | — | — | — | — | 23,300 | 31,100 J | | |
| Manganese | — | — | — | — | — | 6.8 | 3.2 J | | |
| Mercury | — | 0.1 U | 0.1 U | 0.1 U | 0.1 | 0.1 | 0.1 J | | |
| Nickel | — | 1.9 B | 0.5 U | 1.9 B | 0.7 | 0.7 | 1.1 J | | |
| Potassium | — | — | — | — | — | 3,710 | 1,900 | | |
| Selenium | — | 4.0 U | 4.0 R | 4.0 U | 4.4 UJ | 4.4 R | 4.4 R | | |
| Silver | — | 3.3 B | 0.5 U | 3.3 B | 0.4 | 0.4 | 0.6 | | |
| Sodium | — | — | — | — | — | 32,800 | 61,200 | | |
| Thallium | — | 3.6 UJ | 3.6 U | 3.6 UJ | 6.0 J | 2.6 | 2.6 | | |
| Vanadium | — | — | — | — | — | 0.8 | 2.2 J | | |
| Zinc | — | 12.6 B | 6.0 J | 12.6 B | 8.7 | 0.6 UJ | 0.6 | | |
| <u>Volatile Organic Compounds (VOCs)</u> | BRL | BRL | BRL | BRL | BRL | BRL | BRL | | |
| <u>Semi-Volatile Organic Compounds (SVOCs)</u> | BRL | BRL | BRL | BRL | BRL | BRL | BRL | | |
| <u>Pesticides / PCBs</u> | BRL | BRL | BRL | BRL | BRL | BRL | BRL | | |

- Notes:
- 1) All results expressed in micrograms per liter (µg/L).
 - 2) Standard Inorganic Data Qualifiers have been used.
 - 3) Yellow shading indicate a detection above the Contract Required Quantitation Limit (CRQL).
 - 4) Bold red letters with a thick outline indicates a detection above the Trigger Level.
 - 5) BRL = Below Report Limit; reported data values have a data qualifier of U, J, or UJ
 - 6) — = Constituent not analyzed.
 - 7) U = Not detected at the listed reporting limit.
 - 8) B = An estimated value above the method detection limit (MDL) or the instrument detection limit (IDL) but below the CRQL.
 - 9) UJ = A value less than the CRQL but greater than the MDL.
 - 10) J = The analyte was positively identified; the associated numerical value is the approximate concentration of analyte in the sample.
 - 11) R = The sample results are rejected due to deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte can not be verified.
 - 12) CRQL = Contract Required Quantitation Limit
 - 13) Samples analyzed for Dissolved Inorganics were field filtered using a .45 micron, gravity flow filter.
 - 14) Detailed summary tables which list report limits and qualified data values for each compound analyzed for by the laboratory as well as qualified laboratory reports are available upon request.

Skinner Landfill
West Chester, Ohio
Groundwater Analysis Summary Table for Outfall Surface Water Run Off Location SWD-1

| | Sampling Event (All Results Expressed in Units of µg/l) | | | | | | | | |
|---|---|-------------|-----------------|--------|-------------------|-----------------|-----------------|---------------|------|
| | Baseline Results | | | | Quarterly Results | | | | |
| Compound | September-02 | December-02 | February-03 | May-03 | August-03 | November-03 | March-04 | TRIGGER LEVEL | CRQL |
| Inorganics - Metals (Dissolved) ¹³ | | | Location is Dry | | | Location is Dry | Location is Dry | | |
| Antimony | 6.9 B | 3.3 B | — | 5.7 B | 3.7 | — | — | 60 | 60 |
| Arsenic | 3.6 U | 4.3 B | — | 2.9 U | 2.9 | — | — | 10 | 10 |
| Barium | 29.1 B | 33.0 B | — | 22.0 B | 20.8 | — | — | 1,000 | 200 |
| Beryllium | 0.1 U | 0.1 U | — | 0.1 U | 0.1 | — | — | 5 | 5 |
| Cadmium | 0.2 U | 0.2 U | — | 0.2 U | 0.2 | — | — | 5 | 5 |
| Chromium | 0.7 U | 2.1 B | — | 2 B | 0.8 | — | — | 11 | 10 |
| Copper | 2.5 B | 2.4 B | — | 9 B | 5.1 | — | — | 25 | 25 |
| Iron | 1.3 U | 4.9 U | — | 14.1 U | 14.1 | — | — | 5,000 | 100 |
| Lead | 0.1 U | 1.3 U | — | 1.5 UJ | 1.5 | — | — | 4.2 | 3 |
| Mercury | 0.1 U | 0.1 U | — | 0.1 U | 0.1 | — | — | 0.2 | 0.2 |
| Nickel | 2.6 B | 1.7 B | — | 1.3 B | 1.5 | — | — | 96 | 40 |
| Selenium | 4.0 R | 4.0 U | — | 4.4 U | 4.4 | — | — | 5 | 5 |
| Silver | 0.5 U | 2.6 B | — | 0.4 U | 0.4 | — | — | 10 | 10 |
| Thallium | 3.6 U | 3.6 U | — | 2.6 UJ | 2.6 | — | — | 40 | 10 |
| Zinc | 86.6 J | 93.7 | — | 76.2 | 48.0 J | — | — | 86 | 20 |
| Inorganics - Metals and Cyanide (Total) | | | | | | | | | |
| Antimony | — | 6.7 B | — | 5.8 B | 3.7 | — | — | | |
| Arsenic | — | 7.2 B | — | 2.9 U | 2.9 | — | — | | |
| Barium | — | 36.1 B | — | 26.4 B | 21.7 | — | — | | |
| Beryllium | — | 0.1 U | — | 0.1 U | 0.1 | — | — | | |
| Cadmium | — | 0.2 U | — | 0.1 U | 0.2 | — | — | | |
| Chromium | — | 1.9 B | — | 0.2 B | 0.8 | — | — | | |
| Copper | — | 6.3 B | — | 3.6 B | 24.0 | — | — | | |
| Cyanide | 4.0 B | 4.0 U | — | 3.0 U | 3.0 | — | — | 10 | 10 |
| Iron | — | 768 | — | 461 | 72.2 | — | — | | |
| Lead | — | 1.3 U | — | 1.5 UJ | 1.5 | — | — | | |
| Mercury | — | 0.1 U | — | 0.1 U | 0.1 | — | — | | |
| Nickel | — | 3.7 B | — | 1.3 B | 1.4 | — | — | | |
| Selenium | — | 4.0 U | — | 4.4 U | 4.4 UJ | — | — | | |
| Silver | — | 2.0 B | — | 0.4 U | 0.4 | — | — | | |
| Thallium | — | 3.6 UJ | — | 2.6 U | 2.6 UJ | — | — | | |
| Zinc | — | 102 | — | 167 | 91.1 J | — | — | | |
| Volatile Organic Compounds (VOCs) | | | | | | | | | |
| | BRL | BRL | — | BRL | BRL | — | — | | |
| Semi-Volatile Organic Compounds (SVOCs) | | | | | | | | | |
| | BRL | BRL | — | BRL | BRL | — | — | | |
| Pesticides / PCBs | | | | | | | | | |
| | BRL | BRL | — | BRL | BRL | — | — | | |

Notes:

- 1) All results expressed in micrograms per liter (µg/l).
- 2) Standard Inorganic Data Qualifiers have been used.
- 3) Yellow shading indicate a detection above the Contract Required Quantitation Limit (CRQL).
- 4) Bold red letters with a thick outline indicates a detection above the Trigger Level.
- 5) BRL = Below Report Limit; reported data values have a data qualifier of U, J, or UJ
- 6) — = No Sample Available (location is dry)
- 7) U = Not detected at the listed reporting limit.
- 8) B = An estimated value above the method detection limit (MDL) or the instrument detection limit (IDL) but below the CRQL.
- 9) UJ = A value less than the CRQL but greater than the MDL.
- 10) J = The analyte was positively identified; the associated numerical value is the approximate concentration of analyte in the sample.
- 11) R = The sample results are rejected due to deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte can not be verified.
- 12) CRQL = Contract Required Quantitation Limit
- 13) Samples analyzed for Dissolved Inorganics were field filtered using a .45 micron, gravity flow filter.
- 14) Detailed summary tables which list report limits and qualified data values for each compound analyzed for by the laboratory as well as qualified laboratory reports are available upon request.

Skinner Landfill
West Chester, Ohio
Groundwater Analysis Summary Table for Outfall Surface Water Run Off Location SWD-2

| Sampling Event (All Results Expressed in Units of µg/l) | | | | | | | | | |
|---|------------------|-----------------|-------------|--------|-------------------|-----------------|-----------------|---------------|------|
| Compound | Baseline Results | | | | Quarterly Results | | | TRIGGER LEVEL | CRQL |
| | September-02 | December-02 | February-03 | May-03 | August-03 | November-03 | March-04 | | |
| Inorganics - Metals (Dissolved)¹³ | Location is Dry | Location is Dry | | | | Location is Dry | Location is Dry | | |
| Antimony | — | — | 3.9 B | 4.6 B | 3.7 | — | — | 60 | 60 |
| Arsenic | — | — | 3.6 U | 2.9 U | 2.9 | — | — | 20 | 10 |
| Barium | — | — | 15.9 B | 20.4 B | 29.1 | — | — | 1,000 | 200 |
| Beryllium | — | — | 0.1 B | 0.1 U | 0.1 | — | — | 5 | 5 |
| Cadmium | — | — | 0.2 U | 0.2 U | 0.2 | — | — | 5 | 5 |
| Chromium | — | — | 0.7 U | 2.4 B | 0.8 | — | — | 11 | 10 |
| Copper | — | — | 2.3 B | 5.0 B | 1.2 | — | — | 25 | 25 |
| Iron | — | — | 4.9 U | 14.1 U | 14.1 | — | — | 7,000 | 100 |
| Lead | — | — | 1.3 U | 1.5 UJ | 1.5 | — | — | 4.2 | 3 |
| Mercury | — | — | 0.1 U | 0.1 U | 0.1 | — | — | 0.2 | 0.2 |
| Nickel | — | — | 1.8 B | 2.8 B | 2.3 | — | — | 96 | 40 |
| Selenium | — | — | 4.0 U | 4.4 U | 4.4 | — | — | 8.5 | 5 |
| Silver | — | — | 0.5 U | 0.4 U | 0.4 | — | — | 10 | 10 |
| Thallium | — | — | 7.1 B | 2.6 U | 2.6 | — | — | 40 | 10 |
| Zinc | — | — | 1.0 B | 5.5 B | 23.6 J | — | — | 86 | 20 |
| Inorganics - Metals and Cyanide (Total) | | | | | | | | | |
| Antimony | — | — | 8.5 B | 5.0 B | 4.0 | — | — | | |
| Arsenic | — | — | 3.6 U | 2.9 U | 2.9 | — | — | | |
| Barium | — | — | 15.6 B | 21.7 B | 28.4 | — | — | | |
| Beryllium | — | — | 0.1 U | 0.1 U | 0.1 | — | — | | |
| Cadmium | — | — | 0.2 U | 0.2 U | 0.2 | — | — | | |
| Chromium | — | — | 2.2 B | 2.1 B | 0.8 | — | — | | |
| Copper | — | — | 10.9 B | 2.8 B | 11.7 | — | — | | |
| Cyanide | — | — | 4.0 U | — | 3.0 | — | — | 10 | 10 |
| Iron | — | — | 20.6 B | 131 | 17.4 | — | — | | |
| Lead | — | — | 1.3 U | 1.5 UJ | 1.5 | — | — | | |
| Mercury | — | — | 0.1 U | 0.1 U | 0.1 | — | — | | |
| Nickel | — | — | 2.8 B | 2.6 B | 2.4 | — | — | | |
| Selenium | — | — | 4.0 UJ | 4.4 U | 4.4 UJ | — | — | | |
| Silver | — | — | 0.5 U | 0.4 U | 0.4 | — | — | | |
| Thallium | — | — | 3.6 U | 2.6 UJ | 2.6 UJ | — | — | | |
| Zinc | — | — | 4.5 B | 14.7 B | 30.0 J | — | — | | |
| Volatile Organic Compounds (VOCs) | | | | | | | | | |
| bis(2-Chloroethyl) ether | — | — | — | — | — | — | — | 13.6 | 10.0 |
| Semi-Volatile Organic Compounds (SVOCs) | | | | | | | | | |
| Pesticides / PCBs | — | — | BRL | BRL | BRL | — | — | | |

Notes:

Laboratory Analytical data for July, August, and September of 2001 for SWD-2 is labeled as sample location "SWD-4" which was an alternate sample location in very close proximity to SWD-2 used when sample was not available at

- 1) All results expressed in micrograms per liter (µg/L).
- 2) Standard Inorganic Data Qualifiers have been used.
- 3) Yellow shading indicate a detection above the Contract Required Quantitation Limit (CRQL).
- 4) Bold red letters with a thick outline indicates a detection above the Trigger Level.
- 5) BRL = Below Report Limit; reported data values have a data qualifier of U, J, or UJ
- 6) — = No Sample Available (location is dry)
- 7) U = Not detected at the listed reporting limit.
- 8) B = An estimated value above the method detection limit (MDL) or the instrument detection limit (IDL) but below the CRQL.
- 9) UJ = A value less than the CRQL but greater than the MDL.
- 10) J = The analyte was positively identified; the associated numerical value is the approximate concentration of analyte in the sample.
- 11) R = The sample results are rejected due to deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte can not be verified.
- 12) CRQL = Contract Required Quantitation Limit
- 13) Samples analyzed for Dissolved Inorganics were field filtered using a .45 micron, gravity flow filter.
- 14) Detailed summary tables which list report limits and qualified data values for each compound analyzed for by the laboratory as well as qualified laboratory reports are available upon request.

Skinner Landfill
West Chester, Ohio
Groundwater Analysis Summary Table for Outfall Surface Water Run Off Location SWD-3

| Sampling Event (All Results Expressed in Units of µg/l) | | | | | | | | | |
|---|------------------|-------------|-------------|--------|-------------------|-------------|----------|---------------|-------|
| Compound | Baseline Results | | | | Quarterly Results | | | TRIGGER LEVEL | CRQL |
| | September-02 | December-02 | February-03 | May-03 | August-03 | November-03 | March-04 | | |
| Inorganics - Metals (Dissolved) ¹³ | | | | | | | | | |
| Aluminum | — | — | — | — | — | 98.7 | 34.6 | | 200 |
| Antimony | 5.0 B | 6.5 B | 3.5 B | 3.7 U | 3.7 | 3.7 | 3.7 | 60 | 60 |
| Arsenic | 4.2 B | 3.6 U | 3.6 U | 2.9 U | 2.9 | 2.9 | 5.3 | 20 | 10 |
| Barium | 23.4 B | 13.3 B | 34.1 B | 26.6 B | 19.1 | 40.1 | 29.8 | 1,000 | 200 |
| Beryllium | 0.2 B | 0.1 U | 0.3 B | 0.1 U | 0.1 | 0.1 | 0.2 | 5 | 5 |
| Cadmium | 0.2 U | 0.2 U | 0.2 U | 0.2 U | 0.2 | 0.2 | 0.2 | 5 | 5 |
| Calcium | — | — | — | — | — | 130,000 | 125,000 | | 5,000 |
| Chromium | 0.7 U | 2.7 B | 0.7 U | 2.3 B | 0.8 | 1.4 | 0.8 | 11 | 10 |
| Cobalt | — | — | — | — | — | 0.4 | 0.4 | | 50 |
| Copper | 2.0 B | 1.0 U | 1.0 U | 2.3 B | 1.2 | 10.4 | 4.6 J | 25 | 25 |
| Iron | 14.3 B | 4.9 U | 59.5 B | 14.7 B | 14.1 | 59.0 | 17.2 | 7,000 | 100 |
| Lead | 1.3 U | 1.3 U | 1.3 U | 1.5 UJ | 1.5 | 1.5 | 1.5 | 4.2 | 3 |
| Magnesium | — | — | — | — | — | 28,500 | 30,400 | | 5,000 |
| Manganese | — | — | — | — | — | 10.9 | 3.0 | | 15 |
| Mercury | 0.1 U | 0.1 U | 0.1 U | 0.1 U | 0.1 | 0.1 UJ | 0.1 | 0.2 | 0.2 |
| Nickel | 3.7 B | 2.6 B | 0.5 U | 1.1 B | 1.1 | 0.7 | 1.4 | 96 | 40 |
| Potassium | — | — | — | — | — | 3,870 | 3,570 | | 5,000 |
| Selenium | 4.0 U | 4.0 U | 4.0 R | 4.4 U | 4.4 | 4.4 | 4.4 UJ | 8.5 | 5 |
| Silver | 0.5 U | 2.4 B | 0.6 B | 0.4 U | 0.4 | 0.4 | 0.4 | 10 | 10 |
| Sodium | — | — | — | — | — | 11,100 | 12,200 | | 5,000 |
| Thallium | 3.6 U | 3.6 U | 6.5 J | 2.6 UJ | 2.6 | 2.6 | 2.6 | 40 | 10 |
| Vanadium | — | — | — | — | — | 2.2 | 0.8 | | 50 |
| Zinc | 6.7 B | 14.9 B | 3.2 B | 2.2 B | 19.0 J | 91.6 J | 0.6 UJ | 86 | 20 |
| Inorganics - Metals and Cyanide | | | | | | | | | |
| (Total) | | | | | | | | | |
| Aluminum | — | — | — | — | — | 177 | 1,800 | | |
| Antimony | — | 6.5 B | 4.0 B | 4.1 B | 3.7 | 3.7 | 5.2 | | |
| Arsenic | — | 3.6 U | 3.6 U | 2.9 U | 2.9 | 2.9 | 2.9 | | |
| Barium | — | 24.3 B | 41.8 B | 37.7 B | 29.0 | 37.0 | 40.0 | | |
| Beryllium | — | 0.1 U | 0.1 U | 0.1 U | 0.1 | 0.1 | 0.2 | | |
| Cadmium | — | 0.2 U | 0.2 U | 0.2 U | 0.2 | 0.2 | 0.2 | | |
| Calcium | — | — | — | — | — | 12,100 | 131,000 | | |
| Chromium | — | 5.0 B | 6.4 B | 4.2 B | 0.8 | 1.0 | 1.4 | | |
| Cobalt | — | — | — | — | — | 0.4 | 1.5 | | |
| Copper | — | 4.4 B | 8.8 B | 6.0 B | 8.6 | 14.8 | 11.0 J | | |
| Cyanide | 4.0 B | 4.0 U | 4.0 U | 3.0 U | 3.0 | 3.0 | 0.8 | 10 | 10 |
| Iron | — | 3,380 | 1,620 | 3,290 | 3,360 | 155 | 2,200 | | |
| Lead | — | 1.3 U | 1.4 B | 1.5 UJ | 1.5 | 1.5 | 1.5 | | |
| Magnesium | — | — | — | — | — | 26,600 | 31,600 | | |
| Manganese | — | — | — | — | — | 16.5 | 87.5 | | |
| Mercury | — | 0.1 U | 0.1 U | 0.1 U | 0.1 | 0.1 UJ | 0.1 | | |
| Nickel | — | 5.1 B | 2.5 B | 3.4 B | 4.2 | 0.7 | 2.5 | | |
| Potassium | — | — | — | — | — | 3,560 | 4,170 | | |
| Selenium | — | 4.0 U | 4.0 UJ | 4.4 U | 4.4 UJ | 4.4 | 4.4 R | | |
| Silver | — | 2.6 B | 0.5 U | 0.4 U | 0.4 | 0.4 | 0.4 | | |
| Sodium | — | — | — | — | — | 10,300 | 12,600 | | |
| Thallium | — | 3.6 UJ | 3.6 U | 2.6 UJ | 2.6 UJ | 2.6 | 2.6 | | |
| Vanadium | — | — | — | — | — | 0.8 | 2.2 | | |
| Zinc | — | 35.4 | 59.0 | 32.3 | 42.9 J | 32.6 J | 14.6 J | | |
| Volatile Organic Compounds (VOCs) | BRL | BRL | BRL | BRL | BRL | BRL | BRL | | |
| Semi-Volatile Organic Compounds (SVOCs) | BRL | BRL | BRL | BRL | BRL | BRL | BRL | | |
| Pesticides / PCBs | BRL | BRL | BRL | BRL | BRL | BRL | BRL | | |

Notes:

- 1) All results expressed in micrograms per liter (µg/L).
- 2) Standard Inorganic Data Qualifiers have been used.
- 3) Yellow shading indicate a detection above the Contract Required Quantitation Limit (CRQL).
- 4) Bold red letters with a thick outline indicates a detection above the Trigger Level.
- 5) BRL = Below Report Limit; reported data values have a data qualifier of U, J, or UJ
- 6) — = Constituent not analyzed.
- 7) U = Not detected at the listed reporting limit.
- 8) B = An estimated value above the method detection limit (MDL) or the instrument detection limit (IDL) but below the CRQL.
- 9) UJ = A value less than the CRQL but greater than the MDL.
- 10) J = The analyte was positively identified; the associated numerical value is the approximate concentration of analyte in the sample.
- 11) R = The sample results are rejected due to deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte can not be verified.
- 12) CRQL = Contract Required Quantitation Limit
- 13) Samples analyzed for Dissolved Inorganics were field filtered using a .45 micron, gravity flow filter.
- 14) Detailed summary tables which list report limits and qualified data values for each compound analyzed for by the laboratory as well as qualified laboratory reports are available upon request.



LABORATORY DATA VALIDATION REPORT

APPENDIX C

DATA VALIDATION REPORT
FOR
SKINNER LANDFILL SITE
EARTH TECH: PROJECT NUMBER 54280
LABORATORY REPORT NUMBER 204032408
PROJECT MANAGER: Ron Rolker
Date: May 25, 2004
Data Validator: Mark Kromis

APPENDIX C LIST OF ACRONYMS

| | |
|---------------|---|
| BFB | Bromofluorebenzene |
| CC | Continuing Calibration |
| CCV | Continuing Calibration Verification |
| CCB | Continuing Calibration Blanks |
| CLP | Contract Laboratory Program |
| CRDL | Contract Required Detection Limit |
| DFTPP | Decafluorotriphenylphosphine |
| GC/MS | Gas Chromatograph Mass Spectrometer |
| IC | Initial Calibration |
| ICB | Initial Calibration Blank |
| IDL | Instrument Detection Limit |
| ICP | Inductively Coupled Plasma |
| ICS | Interference Check Sample |
| ICV | Initial Calibration Verification |
| ILM | Inorganic Analysis Multi-Media Multi-Concentration |
| INDAM | Individual A Mixture |
| INDBM | Individual B Mixture |
| mg/L | milligrams per liter |
| MS/MSD | Matrix Spike Matrix Spike Duplicate |
| OLC | Organic Analysis Low Concentration |
| OLM | Organic Analysis Multi-Media Multi-Concentration |
| %D | Percent Difference |
| % RSD | Percent Relative Standard Deviation |
| PB | Preparation Blanks |
| QC | Quality Control |
| RF | Response Factor |
| RPD | Relative Percent Difference |
| RRF | Relative Response Factor |
| SDG | Sample Delivery Group |
| SOW | Statement of Work |
| µg/L | micrograms per liter |
| US EPA | United States Environmental Protection Agency |
| VOC | Volatile Organic Compounds |
| VTSR | Validated Time of Sample Receipt |

DATA VALIDATION SUMMARY – SAMPLE DELIVERY GROUP 204032408 INORGANICS

Validation of the inorganics data, as prepared by Gulf Coast Analytical Laboratories (GCAL) for the samples collected from the Skinner Landfill site in March 2004, was conducted by Earth Tech using the National Functional Guidelines for Inorganic Data Review, (US EPA, February, 1994), as appropriate. The results were reported by GCAL under Sample Delivery Group (SDG) 204032408.

| GCAL # | Sample Description |
|-------------|----------------------|
| 20403240801 | SKSWEB1009 |
| 20403240802 | SKSW511009 |
| 20403240803 | SKSW51MS1009 |
| 20403240805 | SKSW51DUP1009 |
| 20403240806 | SKSW521009 |
| 20403240807 | SKSW531009 |
| 20403240808 | SKSWEB1009 (DISS) |
| 20403240809 | SKSW511009 (DISS) |
| 20403240810 | SKSW51MS1009 (DISS) |
| 20403240811 | SKSW51DUP1009 (DISS) |
| 20403240812 | SKSW521009 (DISS) |
| 20403240813 | SKSW531009 (DISS) |

INTRODUCTION

Analyses of metals were performed according to Contract Laboratory Program (CLP)-Inorganic Analysis Multi-media Multi-concentration ILM04.1 Statement of Work (SOW). Results of the sample analyses are reported by the laboratory as either qualified or unqualified. Unqualified results mean that the reported values maybe used without reservation. Various qualifier codes are used by the laboratory to denote specific information regarding the analytical results.

The data validation process is intended to evaluate the data on a technical basis. The data package also was subjected to an internal laboratory quality review prior to submission to Earth Tech for data validation.

During the validation process, laboratory-qualified and unqualified data are verified against all available supporting documentation. Based on this evaluation, qualifier codes may be added, deleted or modified by the data user. Final results are therefore, either qualified or unqualified. Validator-qualified results are annotated with the following codes in accordance with the Functional Guidelines:

- U The constituent was analyzed for, but was not detected above the level of the associated analytical reporting limit. The associated value is either the sample quantitation limit or the sample detection limit.

- J** The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample.
- UJ** The analyte was not detected above the reported sample quantitation limit. However, the reported quantitation limit is approximate and may or may not represent the actual limit of quantitation necessary to accurately and precisely measure the analyte in the sample.
- R** The sample results are rejected due to serious deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte cannot be verified.

Details of the inorganics data validation findings and conclusions are provided in the following sections of this report:

- 1. Holding Times**
- 2. Calibration**
 - A. Initial Calibration (IC)**
 - B. Continuing Calibration (CC)**
- 3. Blanks**
- 4. Inductively Coupled Plasma (ICP) Interference Check Sample**
- 5. Laboratory Control Sample (LCS)**
- 6. Duplicate Analysis**
- 7. Spike Sample Analysis**
- 8. ICP Serial Dilution**
- 9. System Performance**
- 10. Documentation**
- 11. Overall Assessment**

1. HOLDING TIMES

All samples for inorganics analyses were analyzed within the 180-day holding time for preserved aqueous samples. Mercury analyses were conducted within the 28-day holding time for aqueous samples undergoing CLP protocol. Cyanide analyses were conducted within the 14-day holding time. The cooler temperature upon receipt at the laboratory was within the recommended temperature of 4°C +/- 2°C.

2. CALIBRATION

A. Initial Calibration

The percent recoveries for the Initial Calibration Verification (ICV) standard were within Quality Control (QC) limits for all constituents.

B. Continuing Calibration

The percent recoveries for the Continuing Calibration Verification (CCV) standard were within QC limits for all constituents.

3. BLANKS

The Initial Calibration Blank (ICB), Continuing Calibration Blanks (CCB) and Preparation Blanks (PB) were analyzed at the appropriate frequencies. No constituents were detected in the ICB, CCB, and PB blanks above the corresponding Contract Required Detection Limit (CRDL).

4. ICP INTERFERENCE CHECK SAMPLE

Results for the ICP analysis of the Interference Check Sample (ICS) solution AB were within 20% of the true value.

5. LABORATORY CONTROL SAMPLES

Recoveries were within the control limit (80-120%) for all constituents.

6. DUPLICATE ANALYSIS

The laboratory used sample SKSW511009 for the duplicate sample. The Relative Percent Difference (RPD) between the sample and duplicate results for the dissolved fraction were within the acceptance criteria (<20%) for all target compounds. The Relative Percent Difference (RPD) between the sample and duplicate results for the total fraction were within the acceptance criteria (<20%) for all target compounds with the exception of Arsenic, Barium, Chromium, Copper, Iron, Lead, Nickel, and Zinc. As per the National Functional Guidelines, if the results from a duplicate analysis for a particular analyte falls outside the appropriate fixed control windows; qualify the results for that analyte in all associated samples of the same matrix as estimated (J).

7. SPIKE SAMPLE ANALYSIS

The laboratory used sample SKSW511009 for the matrix spike sample. The MS percent recoveries were within the acceptance criteria (75%-125%) with the exception of Antimony (35%), Arsenic (151%), Selenium (0%), and Zinc (128%) in the total fraction and Arsenic (128%) and Selenium (0%) in the dissolved fraction. As per the National Functional Guidelines: if the percent recovery is greater than 30% and less than 74% qualify detected results for that analyte with "J" and non-detected results with "UJ". If the percent recovery is less than 30% qualify detected results for that analyte with "J" and non-detected results with "R".

8. ICP SERIAL DILUTION

As noted in the National Functional Guidelines: If the analyte concentration is at least 50 times above the IDL, its serial dilution analysis must then agree within 10% of the original determination after corrected for dilution. The serial dilution is performed to determine whether any significant chemical or physical interference's exist due to matrix effects. The serial dilution percent differences were within the acceptance criteria for all target analytes with the exception of Barium, Chromium, Iron and Nickel in the total fraction. As per the National Functional Guidelines: if the required 10% Difference criteria are not met, qualify the associated data as estimated "J".

9. SYSTEM PERFORMANCE

The analytical system appears to have been working well at the time of these analyses, based on the evaluation of the raw data.

10. DOCUMENTATION

The documentation appeared accurate and in order.

11. OVERALL ASSESSMENT

The results are acceptable with the validator-added qualifiers.

**DATA VALIDATION SUMMARY – SAMPLE DELIVERY GROUP 204032408
SEMIVOLATILE ORGANICS**

Validation of the Gas Chromatograph/Mass Spectrometer (GC/MS) semi-volatile organics data, as prepared by Gulf Coast Analytical Laboratories (GCAL) for the samples collected from the Skinner Landfill site in March 2004, was conducted by Earth Tech using the National Functional Guidelines for Organic Data Review, (US EPA, October, 1999) as appropriate. The results were reported by GCAL under SDG 204032408.

| GCAL # | Sample Description |
|-------------|--------------------|
| 20403240801 | SKSWEB1009 |
| 20403240802 | SKSW511009 |
| 20403240803 | SKSW51MS1009 |
| 20403240804 | SKSW51MSD1009 |
| 20403240806 | SKSW521009 |
| 20403240807 | SKSW531009 |

INTRODUCTION

Analyses were performed according to CLP-Organic Analysis Multi-Media, Multi-Concentration OLM04.2 SOW. Results of the sample analyses are reported by the laboratory as either qualified or unqualified. Unqualified results mean that the reported values may be used without reservation. Various data qualifier codes are used by the laboratory to denote specific information regarding the analytical results. The data validation process is intended to evaluate the data on a technical basis. The data package also was subjected to an internal laboratory quality review prior to submission to Earth Tech for data validation.

During the validation process, laboratory-qualified and unqualified data are verified against all available supporting documentation. Based on this evaluation, qualifier codes may be added, deleted or modified by the data user. Final results are therefore, either qualified or unqualified. Validator-qualified results are annotated with the following codes in accordance with the Functional Guidelines:

- U The constituent was analyzed for, but was not detected above the level of the associated analytical reporting limit. The associated value is either the sample quantitation limit or the sample detection limit.
- J The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample.
- UJ The analyte was not detected above the reported sample quantitation limit. However, the reported quantitation limit is approximate and may or may not represent the actual limit of quantitation necessary to accurately and precisely measure the analyte in the sample.

- R The sample results are rejected due to serious deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte cannot be verified.

Details of the semivolatile data validation findings and conclusions are provided in the following sections of this report:

1. Holding Times
2. GC/MS Tuning
3. Calibration
 - A. IC
 - B. CC
4. Blanks
5. System Monitoring Compound Recovery
6. MS/MSD
7. Internal Standards Performance
8. Compound Identification
9. Constituent Quantitation and Reported Detection Limits
10. System Performance
11. Documentation
12. Overall Assessment

1. HOLDING TIMES

All samples were initially extracted within the seven-day technical holding time and the five-day VTSR method holding time. The cooler temperature upon receipt at the laboratory was within the recommended temperature of 4°C - - 2°C.

2. GC/MS TUNING

The samples were analyzed on a single GC/MS system, identified as MSSV2. Two decafluorotriphenylphosphine (DFTPP) tunes were run representing the shift in which the standards and samples were analyzed. The DFTPP tunes are acceptable.

3. CALIBRATION

A. Initial Calibration

One IC dated 3/19/04 was analyzed in support of the semivolatile sample analyses. Documentation of the IC was present in the data package, and the Relative Response Factor (RRF), as well as percent % RSD values were accurately reported for all target compounds. The criteria employed for technical data review purposes are different than those used in the method. The laboratory must meet a minimum RRF of 0.01; however, for data review purposes, a RRF criterion of "greater than or equal to 0.05" is applied to all semi-volatile compounds. The RRF's and the average RRF were within the acceptance criteria specified in the method for all reported analytes. The %RSD's were within the acceptance criteria specified in the method for all target analytes with the exception of Di-n-butylphthalate (32.0%), Di-n-octylphthalate (30.3%) and Diethylphthalate (41.8%). The lowest point of the calibration curve was dropped for Di-n-butylphthalate, Di-n-octylphthalate, and Diethylphthalate and the %RSD were recalculated. The recalculated %RSD were within the acceptance criteria of less than 30%. Di-n-octylphthalate and Diethylphthalate were not detected in the associated samples therefore data qualification was not required. The detected results for Di-n-butylphthalate were mitigated do to the presence of Di-n-butylphthalate in the associated method blank.

B. Continuing Calibration

One CC dated 4/7/04 was analyzed in support of the semivolatile sample analyses reported in the data submissions. The RRF's for the CC dated 4/7/04 were within the acceptance criteria. The percent difference (%D) between the average RRF's and the CC Response Factors for the CC dated 4/7/04 were within the acceptance criteria with the exception the %D for Di-n-butylphthalate and Di-n-octylphthalate. As per the National Functional Guidelines, if the %D exceeds the acceptance criteria qualify detected results for that analyte with "J" and non-detected results for that analyte with "UJ".

4. BLANKS

One laboratory semivolatile method blank and an equipment blank were analyzed with this SDG. The results are summarized below.

Method Blank (0325SBLK)

Di-n-butylphthalate was detected at a concentration of 0.851 ppb in method blank 0322SBLK.

Equipment Blank (SKSWEB1009)

Di-n-butylphthalate was detected at a concentration of 0.904 ppb in the equipment blank collected on 3/22/04. The Di-n-butylphthalate result was mitigated by the presence of Di-n-butylphthalate in the associated extraction blank.

5. SYSTEM MONITORING COMPOUND RECOVERY

All reported semivolatile system monitoring compounds were recovered within acceptable control limits.

6. MATRIX SPIKE/MATRIX SPIKE DUPLICATE (MS/MSD)

Sample SKSW511009 was submitted for MS/MSD analysis. The MS/MSD percent recoveries were within the acceptance criteria with the exception of the 4-Nitrophenol. The %RPD between the MS/MSD are within the acceptance criteria. As per the National Functional Guidelines, no action is taken on MS/MSD data alone.

7. INTERNAL STANDARDS PERFORMANCE

Internal standard areas and retention times were within acceptable limits for the reported semivolatile sample analyses.

8. COMPOUND IDENTIFICATION

All reported semivolatile constituents were correctly identified with supporting chromatograms present in the data package.

9. CONSTITUENT QUANTITATION AND REPORTED DETECTION LIMITS

Constituent quantitations were correctly calculated and reported for semivolatile constituents.

10. SYSTEM PERFORMANCE

The analytical system appears to have been working well at the time of these analyses, based on the evaluation of the raw data submitted for review.

11. DOCUMENTATION

The documentation appeared accurate and in order with the exception of the following: GCAL inadvertently left the Date Extracted off of the CLP Form 1's therefore the data validator inserted the Date Extracted on the CLP Form 1's. GCAL also inadvertently left the "B" qualifier off of the CLP Form 1's for the compound Di-n-butylphthalate therefore the data validator inserted a "B" qualifier in the "Q" column of the CLP Form 1's. The "B" qualifier indicates that the analyte was detected in the associated method blank.

12. OVERALL ASSESSMENT

The results are acceptable with the validator-added qualifiers.

DATA VALIDATION SUMMARY – SAMPLE DELIVERY GROUP 204032408 VOLATILE ORGANIC

Validation of the GC/MS volatile organics data, as prepared by Gulf Coast Analytical Laboratories (GCAL) for the samples collected from the Skinner Landfill site in March 2004, was conducted by Earth Tech using the National Functional Guidelines for Organic Data Review, (US EPA, October, 1999), as appropriate. The results were reported by GCAL under SDG 204032408.

| GCAL # | Sample Description |
|-------------|--------------------|
| 20403240801 | SKSWEB1009 |
| 20403240802 | SKSW511009 |
| 20403240803 | SKSW51MS1009 |
| 20403240804 | SKSW51MSD1009 |
| 20403240806 | SKSW521009 |
| 20403240807 | SKSW531009 |

INTRODUCTION

Analyses were performed according to CLP-Organic Analysis Low Concentration OLC02.0 SOW. Results of the sample analyses are reported by the laboratory as either qualified or unqualified. Unqualified results mean that the reported values may be used without reservation. Various qualifier codes are used by the laboratory to denote specific information regarding the analytical results. The data validation process is intended to evaluate the data on a technical basis. The data package also was subjected to an internal laboratory quality review prior to submission to Earth Tech for data validation.

During the validation process, laboratory-qualified and unqualified data are verified against all available supporting documentation. Based on this evaluation, qualifier codes may be added, deleted or modified by the data user. Final results are therefore, either qualified or unqualified. Validator-qualified results are annotated with the following codes in accordance with the Functional Guidelines:

- U The constituent was analyzed for, but was not detected above the level of the associated analytical reporting limit. The associated value is either the sample quantitation limit or the sample detection limit.
- J The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample.
- UJ The analyte was not detected above the reported sample quantitation limit.

However, the reported quantitation limit is approximate and may or may not represent the actual limit of quantitation necessary to accurately and precisely measure the analyte in the sample.

- R** The sample results are rejected due to serious deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte cannot be verified.

The volatiles data validation findings and conclusions are provided in the following sections of this report:

- 1. Holding Times**
- 2. GC/MS Tuning**
- 3. Calibration**
 - A. IC**
 - B. CC**
- 4. Blanks**
- 5. System Monitoring Compound Recovery**
- 6. MS/MSD**
- 7. Laboratory Control Sample**
- 8. Internal Standards Performance**
- 9. Compound Identification**
- 10. Constituent Quantitation and Reported Detection Limits**
- 11. System Performance**
- 12. Documentation**
- 13. Overall Assessment**

1. HOLDING TIMES

All samples for Volatile Organic Compounds (VOC) analyses were analyzed within the 14-day technical holding time and the 10-day VTSR method holding time. The cooler temperature upon receipt at the laboratory was within the recommended temperature of 4°C +/- 2°C.

2. GC/MS TUNING

The samples were analyzed on two GC/MS system, identified as MSV0 and MSV2. Two bromofluorobenzene (BFB) tunes were run. The BFB tunes are acceptable.

3. CALIBRATION

A. Initial Calibration

One IC dated 3/23/04 was analyzed on instrument MSV2 in support of the volatile sample analyses reported in the data submissions. Documentation of the IC standards is present in the data package, and RRF's as well as %RSD values were accurately reported. The criteria employed for technical data review purposes are different than those used in the method. The laboratory must meet a minimum RRF of 0.01; however, for data review purposes, a RRF criterion of "greater than or equal to 0.05" is applied to all volatile compounds. The %RSD's were within the acceptance criteria specified in the method for all target analytes. The RRF's and the average RRF for the IC dated 3/23/04 were within the acceptance criteria specified in the method for all target analytes with the exception of Acetone. As per the National Functional Guidelines, if any initial calibration RRF is less than 0.05, qualify positive results that have acceptable mass spectral identification with "J", using professional judgement, and non-detected analytes as unusable (R). It should be noted that the laboratory did meet the minimum RRF of 0.01 for all target compounds.

B. Continuing Calibration

One CC dated 3/23/04 was analyzed on instrument MSV2 in support of the volatile sample analyses reported in the data submissions. The percent difference (%D) between the average RRF's and the CC RF's were within the acceptance criteria for all target analytes. The CC RRF's were within the acceptance criteria specified in the method for all target analytes with the exception of Acetone. The Acetone results were previously qualified under section 3A above.

4. BLANKS

One laboratory volatile method blank, storage blank, and a Equipment Blank were analyzed with this SDG. The results are summarized below.

MB154405

1,2-Dichlorobenzene was detected at a concentration of 0.17 ppb in the method blank analyzed on 3/24/04.

Storage Blank (VHBLK01)

No compounds were detected above the MDL in the storage blank analyzed on 3/24/04.

Trip Blank

There was no Trip Blank submitted for this sampling event.

Equipment Blank (SKSWEB1009)

Acetone, Methylene chloride and Toluene were detected at concentrations of 2.8 ppb, 0.47 ppb and 1.1 ppb respectively in the equipment blank collected on 3/22/04.

5. SYSTEM MONITORING COMPOUND RECOVERY

All reported volatile system monitoring compounds were recovered within acceptable control limits for all samples.

6. MATRIX SPIKE/MATRIX SPIKE DUPLICATE

Sample SKSW511009 was submitted for MS/MSD analysis. The MS/MSD percent recoveries and %RPD between the MS/MSD were within the acceptance criteria.

7. LABORATORY CONTROL SAMPLE

A LCS was analyzed in conjunction with this SDG. Recoveries were within the control limit for all constituents.

8. INTERNAL STANDARDS PERFORMANCE

Internal Standard (IS) areas and retention times were within acceptable limits for the reported volatile sample analyses.

9. COMPOUND IDENTIFICATION

All reported VOCs were correctly identified with supporting chromatograms present in the data package.

10. CONSTITUENT QUANTITATION AND REPORTED DETECTION LIMITS

Constituent quantitations were correctly calculated and reported for VOCs with the exception of Ethylbenzene. The Ethylbenzene standard and detected results were originally quantitated using the incorrect quantitation ion (GCAL used 106 instead of 91). GCAL corrected the mistake and re-submitted the corrected pages that were affected in the laboratory report. The overall effect had no impact in the final result for Ethylbenzene.

11. SYSTEM PERFORMANCE

The analytical system appears to have been working well at the time of these analyses, based on the evaluation of the raw data.

12. DOCUMENTATION

The documentation appeared accurate and in order with the exception of the following: GCAL inadvertently transposed the area counts and retention times for IS Chlorobenzene-d5 and 1,4-Dichlorobenzene-d4. The data validator corrected the mistake by drawing arrows to indicate the correct area counts and retention times for IS Chlorobenzene-d5 and 1,4-Dichlorobenzene-d4.

13. OVERALL ASSESSMENT

The results are acceptable with the validator-added qualifiers.

DATA VALIDATION SUMMARY - SAMPLE DELIVERY GROUP 204032408 PESTICIDES

Validation of the Gas Chromatography (GC) pesticides data, as prepared by Gulf Coast Analytical Laboratories (GCAL) for the samples collected from the Skinner Landfill site in March 2004, was conducted by Earth Tech using the National Functional Guidelines for Organic Data Review, (US EPA, October, 1999), as appropriate. The results were reported by GCAL under SDG 204032408.

| GCAL # | Sample Description |
|-------------|--------------------|
| 20403240801 | SKSWEB1009 |
| 20403240802 | SKSW511009 |
| 20403240803 | SKSW51MS1009 |
| 20403240804 | SKSW51MSD1009 |
| 20403240806 | SKSW521009 |
| 20403240807 | SKSW531009 |

INTRODUCTION

Analyses were performed according to CLP-Organic Analysis Multi-Media, Multi-Concentration OLM04.2 SOW. Results of the sample analyses are reported by the laboratory as either qualified or unqualified. Unqualified results mean that the reported values may be used without reservation. Various qualifier codes are used by the laboratory to denote specific information regarding the analytical results.

The data validation process is intended to evaluate the data on a technical basis. The data package also was subjected to an internal laboratory quality review prior to submission to Earth Tech for data validation.

During the validation process, laboratory-qualified and unqualified data are verified against all available supporting documentation. Based on this evaluation, qualifier codes may be added, deleted or modified by the data user. Final results are therefore, either qualified or unqualified. Validator-qualified results are annotated with the following codes in accordance with the Functional Guidelines:

- U** The constituent was analyzed for, but was not detected above the level of the associated analytical reporting limit. The associated value is either the sample quantitation limit or the sample detection limit.
- J** The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample.
- UJ** The analyte was not detected above the reported sample quantitation limit. However, the reported quantitation limit is approximate and may or may not represent the actual limit of quantitation necessary to accurately and precisely measure the analyte in the sample

- R The sample results are rejected due to serious deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte cannot be verified.

Details of the pesticide data validation findings and conclusions are provided in the following sections of this report:

1. Holding Times
2. Gas Chromatograph/Electronic Capture Detector (GC/ECD) Instrument Performance Check
3. IC
4. Calibration Verification
5. Blanks
6. Surrogate Spikes
7. Matrix Spike/Matrix Spike Duplicate (MS/MSD)
8. Pesticide Cleanup Checks
9. Target Compound Identification
10. Constituent Quantitation and Reported Detection Limits
11. Documentation
12. Overall Assessment

1. HOLDING TIMES

All samples were extracted within the seven-day technical holding time and the five-day VTSR method holding time. The cooler temperature upon receipt at the laboratory was within the recommended temperature of 4°C +/- 2°C.

2. GC/ECD INSTRUMENT PERFORMANCE CHECK

The Performance Evaluation Mixture (PEM) was analyzed at the correct frequency. Absolute retention times were within limits.

The percent resolution between adjacent peaks was within QC limits for the Pesticide Analyte Resolution Check.

The percent resolution between adjacent peaks is within QC limits for the Performance Evaluation Mixtures (PEM). The percent breakdown for both 4,4-DDT and Endrin in each PEM was less than 20.0% for both GC columns. The combined percent breakdown for 4,4-DDT and Endrin in each PEM was less than 30.0% for both GC columns.

3. INITIAL CALIBRATION

Individual standard mixtures A and B were analyzed at the correct frequencies and concentrations. The percent resolution criterion was met for Individual standard mixtures A and B.

The Percent Relative Standard Deviation (%RSD) of the calibration factors for each of the single component pesticides was less than 20%.

The multi-component target compounds were analyzed separately on both columns at a single concentration level. Retention times were determined from a minimum of three peaks.

4. CALIBRATION VERIFICATION

Absolute retention times were within appropriate time retention windows

5. BLANKS

One laboratory method blank was analyzed with this SDG. The results are summarized below.

Method Blank 155051

No constituents were detected above the laboratory-reporting limit. This blank corresponds to all samples extracted on 3/26/04.

Equipment Blank SKSWEB1009

No constituents were detected above the laboratory-reporting limit in the equipment blank collected on 3/22/04.

6. SURROGATE SPIKES

Decachlorobiphenyl (DCB) and tetrachloro-m-xylene (TCX) surrogate spike recoveries were within the acceptance criteria for all samples.

7. MATRIX SPIKE/MATRIX SPIKE DUPLICATE

Sample SKSW511009 was submitted for MS/MSD analysis. The MS/MSD percent recoveries were within the acceptance criteria. The %RPD between the MS/MSD are within the acceptance criteria.

8. PESTICIDE CLEANUP CHECKS

Recoveries of all pesticides and surrogates were within 80-120% for the lot of Florisil cartridges utilized for pesticide cleanup with the exception of 4,4'-DDT (130%). There were no target compounds detected in the associated samples therefore no action was taken.

9. TARGET COMPOUND IDENTIFICATION

All reported pesticide data were correctly identified with supporting chromatograms present in the data package.

10. CONSTITUENT QUANTITATION AND REPORTED DETECTION LIMITS

Constituent quantitations were correctly calculated and reported for pesticide constituents.

11. DOCUMENTATION

The documentation appeared accurate and in order with the exception of the following: GCAL inadvertently left the Date Extracted off of the CLP Form 1's therefore the data validator inserted the Date Extracted on the CLP Form 1's.

12. OVERALL ASSESSMENT

The results are acceptable with the validator-added qualifiers.

REFERENCES

US EPA, 1994. *National Functional Guidelines for Inorganic Data Review.*

US EPA, 1999. *National Functional Guidelines for Organic Data Review.*



ANALYTICAL RESULTS

PERFORMED BY

GULF COAST ANALYTICAL LABORATORIES, INC.

Report Date 04/16/2004

GCAL Report 204032408

Deliver To Earth Tech
200 Vine Street
Wilder, KY 41076
859-442-2300

Attn Pat Higgins

Customer Earth Tech

Project Skinner Landfill

000001

CASE NARRATIVE

Client: Earth Tech **Report:** 204032408

Gulf Coast Analytical Laboratories received and analyzed the sample(s) listed on the sample cross-reference page of this report. Receipt of the sample(s) is documented by the attached chain of custody. This applies only to the sample(s) listed in this report. No sample integrity or quality control exceptions were identified unless noted below.

SEMI-VOLATILES MASS SPECTROMETRY

The MS/MSD recoveries for 4-Nitrophenol are above the control limits.

SEMI-VOLATILES GAS CHROMATOGRAPHY

In the Florisil check analysis, the recovery for DDT was above recovery limit, however DDT was not detected in the associated samples.

METALS

Barium, Chromium, Iron, and Nickel are flagged as estimated for samples associated with prep batch 271292 due to the fact that the percent difference between the original sample result and the serial dilution result is greater than 10. A chemical or physical interference is suspected. In the ILM04.1 - CLP Metals analysis, the MS recovery was outside the control limits for Antimony, Arsenic, Selenium, and Zinc. The LCS recovery was within control limits. This indicates the analysis is in control and the sample is affected by matrix interference. A post-digestion spike was performed on the QC sample for this batch with a recovery of 109% for Antimony, 131% for Arsenic, 64% for Selenium, and 139% for Zinc. The MS recovery is not applicable for Iron and Lead because the sample concentration is greater than four times the spike concentration. The Sample/Duplicate RPD for Arsenic, Barium, Chromium, Copper, Iron, Lead, Nickel, and Zinc was outside the control limits. The heterogeneous nature of the QC sample is believed to be responsible for this.

In the ILM04.1 - CLP Metals analysis for prep batch 271294, the MS recovery was outside the control limits for Arsenic and Selenium. The LCS recovery was within control limits. This indicates the analysis is in control and the sample is affected by matrix interference. A post-digestion spike was performed on the QC sample for this batch with a recovery of 106% for Arsenic and 94% for Selenium.

Laboratory Endorsement

Sample analysis was performed in accordance with approved methodologies provided by the Environmental Protection Agency or other recognized agencies. The samples and their corresponding extracts will be maintained for a period of 30 days unless otherwise arranged. Following this retention period the samples will be disposed in accordance with GCAL's Standard Operating Procedures.

Common Abbreviations Utilized in this Report

ND Indicates the result was Not Detected at the specified RDL
DO Indicates the result was Diluted Out
MI Indicates the result was subject to Matrix Interference
TNTC Indicates the result was Too Numerous To Count
SUBC Indicates the analysis was Sub-Contracted
FLD Indicates the analysis was performed in the Field
PQL Practical Quantitation Limit
MDL Method Detection Limit
RDL Reporting Detection Limit
00:00 Reported as a time equivalent to 12:00 AM

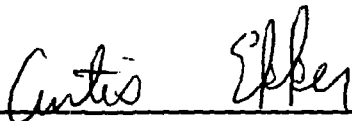
Reporting Flags Utilized in this Report

J Indicates an estimated value
U Indicates the compound was analyzed for but not detected
B (ORGANICS) Indicates the analyte was detected in the associated Method Blank
B (INORGANICS) Indicates the result is between the RDL and MDL

Sample receipt at GCAL is documented through the attached chain of custody. In accordance with ISO Guide 25 and NELAC, this report shall be reproduced only in full and with the written permission of GCAL. The results contained within this report relate only to the samples reported. The documented results are presented within this report.

This report pertains only to the samples listed in the Report Sample Summary and should be retained as a permanent record thereof. The results contained within this report are intended for the use of the client. Any unauthorized use of the information contained in this report is prohibited.

I certify that this data package is in compliance with the terms and conditions of the contract and Statement of Work both technically and for completeness, for other than the conditions in the case narrative. Release of the data contained in this hardcopy data package and in the computer-readable data submitted has been authorized by the Quality Assurance Manager or his/her designee, as verified by the following signature.



CURTIS EKKER
DATA VALIDATION MANAGER
GCAL REPORT 204032408

THIS REPORT CONTAINS 436 PAGES.

000003

Report Sample Summary

| GCAL ID | Client ID | Matrix | Collect Date/Time | Receive Date/Time |
|-------------|----------------------|--------|-------------------|-------------------|
| 20403240801 | SKSWEB1009 | Water | 03/22/2004 13:45 | 03/24/2004 09:20 |
| 20403240802 | SKSW511009 | Water | 03/22/2004 10:40 | 03/24/2004 09:20 |
| 20403240803 | SKSW51MS1009 | Water | 03/22/2004 10:55 | 03/24/2004 09:20 |
| 20403240804 | SKSW51MSD1009 | Water | 03/22/2004 11:15 | 03/24/2004 09:20 |
| 20403240805 | SKSW51DUP1009 | Water | 03/22/2004 11:15 | 03/24/2004 09:20 |
| 20403240806 | SKSW521009 | Water | 03/22/2004 12:30 | 03/24/2004 09:20 |
| 20403240807 | SKSW531009 | Water | 03/22/2004 12:55 | 03/24/2004 09:20 |
| 20403240808 | SKSWEB1009 (DISS) | Water | 03/22/2004 13:45 | 03/24/2004 09:20 |
| 20403240809 | SKSW511009 (DISS) | Water | 03/22/2004 10:40 | 03/24/2004 09:20 |
| 20403240810 | SKSW51MS1009 (DISS) | Water | 03/22/2004 10:55 | 03/24/2004 09:20 |
| 20403240811 | SKSW51DUP1009 (DISS) | Water | 03/22/2004 11:15 | 03/24/2004 09:20 |
| 20403240812 | SKSW521009 (DISS) | Water | 03/22/2004 12:30 | 03/24/2004 09:20 |
| 20403240813 | SKSW531009 (DISS) | Water | 03/22/2004 12:55 | 03/24/2004 09:20 |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SKSWEB1009

Lab Name: GCAL

Contract:

Lab Code: LA024

Case No.:

SAS No.:

SDG No.: 204032408

Matrix: (soil/water) Water

Sample wt/vol: 25 (g/ml) mL

Lab Sample ID: 20403240801

Level: (low/med)

Lab File ID: 2040324/T2586

% Moisture: not dec.

Date Collected: 03/22/04 Time: 1345

GC Column: DB-624-30M ID: .53 (mm)

Date Received: 03/24/04

Instrument ID: MSV2

Date Analyzed: 03/24/04 Time: 1851

Soil Extract Volume: (µL)

Dilution Factor: 1 Analyst: RJO

Soil Aliquot Volume: (µL)

Prep Batch: Analytical Batch: 271209

CONCENTRATION UNITS: ug/L

Analytical Method: OLCO 2.1

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|------------|---------------------------|--------|---|-------|-----|
| 71-55-6 | 1,1,1-Trichloroethane | 1.0 | U | 0.010 | 1.0 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1.0 | U | 0.010 | 1.0 |
| 79-00-5 | 1,1,2-Trichloroethane | 1.0 | U | 0.010 | 1.0 |
| 75-34-3 | 1,1-Dichloroethane | 1.0 | U | 0.010 | 1.0 |
| 75-35-4 | 1,1-Dichloroethene | 1.0 | U | 0.010 | 1.0 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 106-93-4 | 1,2-Dibromoethane | 1.0 | U | 0.010 | 1.0 |
| 95-50-1 | 1,2-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 107-06-2 | 1,2-Dichloroethane | 1.0 | U | 0.010 | 1.0 |
| 540-59-0 | 1,2-Dichloroethene | 1.0 | U | 0.010 | 1.0 |
| 78-87-5 | 1,2-Dichloropropane | 1.0 | U | 0.010 | 1.0 |
| 541-73-1 | 1,3-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 106-46-7 | 1,4-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 78-93-3 | 2-Butanone | 5.0 | U | 0.010 | 5.0 |
| 591-78-6 | 2-Hexanone | 5.0 | U | 0.010 | 5.0 |
| 108-10-1 | 4-Methyl-2-pentanone | 5.0 | U | 0.010 | 5.0 |
| 67-64-1 | Acetone | 2.8 | J | 0.010 | 5.0 |
| 71-43-2 | Benzene | 1.0 | U | 0.010 | 1.0 |
| 75-27-4 | Bromodichloromethane | 1.0 | U | 0.010 | 1.0 |
| 75-25-2 | Bromoform | 1.0 | U | 0.010 | 1.0 |
| 74-83-9 | Bromomethane | 1.0 | U | 0.010 | 1.0 |
| 75-15-0 | Carbon disulfide | 1.0 | U | 0.010 | 1.0 |
| 56-23-5 | Carbon tetrachloride | 1.0 | U | 0.010 | 1.0 |
| 108-90-7 | Chlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 75-00-3 | Chloroethane | 1.0 | U | 0.010 | 1.0 |
| 67-66-3 | Chloroform | 1.0 | U | 0.010 | 1.0 |
| 74-87-3 | Chloromethane | 1.0 | U | 0.010 | 1.0 |
| 124-48-1 | Dibromochloromethane | 1.0 | U | 0.010 | 1.0 |
| 10061-01-5 | cis-1,3-Dichloropropene | 1.0 | U | 0.010 | 1.0 |
| 10061-02-6 | trans-1,3-Dichloropropene | 1.0 | U | 0.010 | 1.0 |
| 100-41-4 | Ethylbenzene | 1.0 | U | 0.010 | 1.0 |
| 75-09-2 | Methylene chloride | 0.47 | J | 0.010 | 2.0 |
| 100-42-5 | Styrene | 1.0 | U | 0.010 | 1.0 |

VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SKSWEB1009

Lab Name: GCAL

Contract:

Lab Code: LA024

Case No:

SAS No:

SDG No.: 204032408

Matrix: (soil/water) Water

Sample wt/vol: 25 (g/mL) mL

Lab Sample ID: 20403240801

Level: (low/med)

Lab File ID: 2040324/T2586

% Moisture: not dec.

Date Collected: 03/22/04

Time: 1345

GC Column: DB-624-30M ID: 53 (mm)

Date Received: 03/24/04

Instrument ID: MSV2

Date Analyzed: 03/24/04

Time: 1851

Soil Extract Volume: (µL)

Dilution Factor: 1

Analyst: RJO

Soil Aliquot Volume: (µL)

Prep Batch:

Analytical Batch: 271209

CONCENTRATION UNITS: µg/L

Analytical Method: OLCO 2.1

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|-----------|-------------------|--------|---|-------|-----|
| 127-18-4 | Tetrachloroethene | ND | U | 0.010 | 1.0 |
| 106-86-3 | Toluene | ND | | 0.010 | 1.0 |
| 79-01-6 | Trichloroethene | ND | U | 0.010 | 1.0 |
| 75-01-4 | Vinyl chloride | ND | U | 0.010 | 1.0 |
| 1330-20-7 | Xylene (total) | ND | U | 0.010 | 1.0 |

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

SKSWEB1009

Lab Name: GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204032408
Matrix: Water Lab Sample ID: 20403240801
Sample wt/vol: _____ Units: _____ Lab File ID: _____
Level: (low/med) _____ Date Collected: 03/22/04 Time: 1345
% Moisture: not dec. _____ Date Received: 03/24/04
GC Column: DB-624-30M ID: .25 (mm) Date Analyzed: 03/24/04 Time: 1851
Instrument ID: MSV2 Dilution Factor: 1 Analyst: RSP
Soil Extract Volume: _____ (µL)
Soil Aliquot Volume: _____ (µL)

Number TICs Found: 0

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | RT | EST. CONC. | Q |
|---------|------------------|----|------------|---|
| 1. | No tics detected | | | |

VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SKSW511009

Lab Name: GCAL

Contract:

Lab Code: LA024

Case No.:

SAS No.:

SDG No.: 204032408

Matrix: (soil/water) Water

Sample wt/vol: 25 (g/ml) mL

Lab Sample ID: 20403240802

Level: (low/med)

Lab File ID: 2040324/T2579

% Moisture: not dec.

Date Collected: 03/22/04

Time: 1040

GC Column: DB-624-30M

ID: 53

(mm)

Date Received: 03/24/04

Instrument ID: MSV2

Date Analyzed: 03/24/04

Time: 1541

Soil Extract Volume: (µL)

Dilution Factor: 1

Analyst: RJO

Soil Aliquot Volume: (µL)

Prep Batch:

Analytical Batch: 271209

CONCENTRATION UNITS µg/L

Analytical Method: OLCO 2.1

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|------------|---------------------------|--------|---|-------|-----|
| 71-55-6 | 1,1,1-Trichloroethane | 1.0 | U | 0.010 | 1.0 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1.0 | U | 0.010 | 1.0 |
| 79-00-5 | 1,1,2-Trichloroethane | 1.0 | U | 0.010 | 1.0 |
| 75-34-3 | 1,1-Dichloroethane | 1.0 | U | 0.010 | 1.0 |
| 75-35-4 | 1,1-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 106-83-4 | 1,2-Dibromoethane | 1.0 | U | 0.010 | 1.0 |
| 95-50-1 | 1,2-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 107-06-2 | 1,2-Dichloroethane | 1.0 | U | 0.010 | 1.0 |
| 540-59-0 | 1,2-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 78-87-5 | 1,2-Dichloropropane | 1.0 | U | 0.010 | 1.0 |
| 541-73-1 | 1,3-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 106-46-7 | 1,4-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 78-83-3 | 2-Butanone | 5.0 | U | 0.010 | 5.0 |
| 581-78-6 | 2-Hexanone | 5.0 | U | 0.010 | 5.0 |
| 106-10-1 | 4-Methyl-2-pentanone | 5.0 | U | 0.010 | 5.0 |
| 67-64-1 | Acetone | 5.0 | U | 0.010 | 5.0 |
| 71-43-2 | Benzene | 1.0 | U | 0.010 | 1.0 |
| 75-27-4 | Bromodichloromethane | 1.0 | U | 0.010 | 1.0 |
| 75-25-2 | Bromoform | 1.0 | U | 0.010 | 1.0 |
| 74-83-9 | Bromomethane | 1.0 | U | 0.010 | 1.0 |
| 75-15-0 | Carbon disulfide | 1.0 | U | 0.010 | 1.0 |
| 56-23-5 | Carbon tetrachloride | 1.0 | U | 0.010 | 1.0 |
| 106-90-7 | Chlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 75-00-3 | Chloroethane | 1.0 | U | 0.010 | 1.0 |
| 67-66-3 | Chloroform | 1.0 | U | 0.010 | 1.0 |
| 74-87-3 | Chloromethane | 1.0 | U | 0.010 | 1.0 |
| 124-48-1 | Dibromochloromethane | 1.0 | U | 0.010 | 1.0 |
| 10061-02-6 | trans-1,3-Dichloropropene | 1.0 | U | 0.010 | 1.0 |
| 10061-01-5 | cis-1,3-Dichloropropene | 1.0 | U | 0.010 | 1.0 |
| 100-41-4 | Ethylbenzene | 1.0 | U | 0.010 | 1.0 |
| 75-09-2 | Methylene chloride | 2.0 | U | 0.010 | 2.0 |
| 100-42-5 | Styrene | 1.0 | U | 0.010 | 1.0 |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SKSW511009

Lab Name: GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204032408
Matrix: (soil/water) Water
Sample wt/vol: 25 (g/ml) mL Lab Sample ID: 20403240802
Level: (low/med) _____ Lab File ID: 2040324/T2579
% Moisture: not dec. _____ Date Collected: 03/22/04 Time: 1040
GC Column: DB-624-30M ID: .53 (mm) Date Received: 03/24/04
Instrument ID: MSV2 Date Analyzed: 03/24/04 Time: 1541
Soil Extract Volume: _____ (µL) Dilution Factor: 1 Analyst: RJO
Soil Aliquot Volume: _____ (µL) Prep Batch: _____ Analytical Batch: 271209
Analytical Method: OLCO 2.1
CONCENTRATION UNITS: ug/L

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|-----------|-------------------|--------|---|-------|-----|
| 127-18-4 | Tetrachloroethene | 1.0 | U | 0.010 | 1.0 |
| 108-88-3 | Toluene | 1.0 | U | 0.010 | 1.0 |
| 79-01-6 | Trichloroethene | 1.0 | U | 0.010 | 1.0 |
| 75-01-4 | Vinyl chloride | 1.0 | U | 0.010 | 1.0 |
| 1330-20-7 | Xylene (total) | 1.0 | U | 0.010 | 1.0 |

1E
VOLATILE ORGANIC ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

SKSW511009

Lab Name: GCAL Contract: _____
 Lab Code: LA024 Case No. _____ SAS No. _____ SDG No.: 204032408
 Matrix: Water Lab Sample ID: 20403240802
 Sample w/vol: _____ Units _____ Lab File ID: _____
 Level: (low/med) _____ Date Collected: 03/22/04 Time: 1040
 % Moisture: not dec. _____ Date Received: 03/24/04
 GC Column: DB-624-30M ID: .25 (mm) Date Analyzed: 03/24/04 Time: 1541
 Instrument ID: MSV2 Dilution Factor: 1 Analyst: RSP
 Soil Extract Volume: _____ (µL)
 Soil Aliquot Volume: _____ (µL)

Number TICs Found: 0

CONCENTRATION UNITS

| CAS NO. | COMPOUND | RT | EST. CONC. | Q |
|---------|------------------|----|------------|---|
| 1. | No tics detected | | | |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SKSW521009

Lab Name: GCAL

Contract:

Lab Code: LA024

Case No.:

SAS No.:

SDG No.: 204032408

Matrix: (soil/water) Water

Sample wt/vol: 25 (g/ml) mL

Lab Sample ID: 20403240806

Level: (low/med)

Lab File ID: 2040324/T2584

% Moisture: not dec.

Date Collected: 03/22/04 Time: 1230

GC Column: DB-624-30M ID: .53 (mm)

Date Received: 03/24/04

Instrument ID: MSV2

Date Analyzed: 03/24/04 Time: 1801

Soil Extract Volume: (µL)

Dilution Factor: 1 Analyst: RJO

Soil Aliquot Volume: (µL)

Prep Batch: Analytical Batch: 271209

CONCENTRATION UNITS: ug/L

Analytical Method: OLCO 2.1

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|------------|---------------------------|--------|---|-------|-----|
| 71-55-6 | 1,1,1-Trichloroethane | 1.0 | U | 0.010 | 1.0 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1.0 | U | 0.010 | 1.0 |
| 79-00-5 | 1,1,2-Trichloroethane | 1.0 | U | 0.010 | 1.0 |
| 75-34-3 | 1,1-Dichloroethane | 1.0 | U | 0.010 | 1.0 |
| 75-35-4 | 1,1-Dichloroethene | 1.0 | U | 0.010 | 1.0 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 106-93-4 | 1,2-Dibromoethane | 1.0 | U | 0.010 | 1.0 |
| 95-50-1 | 1,2-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 107-06-2 | 1,2-Dichloroethane | 1.0 | U | 0.010 | 1.0 |
| 540-59-0 | 1,2-Dichloroethene | 1.0 | U | 0.010 | 1.0 |
| 78-87-5 | 1,2-Dichloropropane | 1.0 | U | 0.010 | 1.0 |
| 541-73-1 | 1,3-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 106-46-7 | 1,4-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 78-93-3 | 2-Butanone | 5.0 | U | 0.010 | 5.0 |
| 591-78-6 | 2-Hexanone | 5.0 | U | 0.010 | 5.0 |
| 108-10-1 | 4-Methyl-2-pentanone | 5.0 | U | 0.010 | 5.0 |
| 67-64-1 | Acetone | 5.0 | U | 0.010 | 5.0 |
| 71-43-2 | Benzene | 1.0 | U | 0.010 | 1.0 |
| 75-27-4 | Bromodichloromethane | 1.0 | U | 0.010 | 1.0 |
| 75-25-2 | Bromoform | 1.0 | U | 0.010 | 1.0 |
| 74-83-9 | Bromomethane | 1.0 | U | 0.010 | 1.0 |
| 75-15-0 | Carbon disulfide | 1.0 | U | 0.010 | 1.0 |
| 56-23-5 | Carbon tetrachloride | 1.0 | U | 0.010 | 1.0 |
| 108-90-7 | Chlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 75-00-3 | Chloroethane | 1.0 | U | 0.010 | 1.0 |
| 67-66-3 | Chloroform | 1.0 | U | 0.010 | 1.0 |
| 74-87-3 | Chloromethane | 1.0 | U | 0.010 | 1.0 |
| 124-48-1 | Dibromochloromethane | 1.0 | U | 0.010 | 1.0 |
| 10061-01-5 | cis-1,3-Dichloropropene | 1.0 | U | 0.010 | 1.0 |
| 10061-02-6 | trans-1,3-Dichloropropene | 1.0 | U | 0.010 | 1.0 |
| 100-41-4 | Ethylbenzene | 1.0 | U | 0.010 | 1.0 |
| 75-09-2 | Methylene chloride | 2.0 | U | 0.010 | 2.0 |
| 100-42-5 | Styrene | 1.0 | U | 0.010 | 1.0 |

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VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SKSW521009

Lab Name: GCAL

Contract: _____

Lab Code: LA024

Case No.: _____

SAS No. _____

SDG No.: 204032408

Matrix: (soil/water) Water

Sample w/vol: 25 (g/ml) mL

Lab Sample ID: 20403240806

Level: (low/med) _____

Lab File ID: 2040324/T2584

% Moisture: not dec. _____

Date Collected: 03/22/04

Time: 1230

GC Column: DB-624-30M ID: 53 (mm)

Date Received: 03/24/04

Instrument ID: MSV2

Date Analyzed: 03/24/04

Time: 1801

Soil Extract Volume: _____ (μL)

Dilution Factor: 1

Analyst: RJO

Soil Aliquot Volume: _____ (μL)

Prep Batch: _____

Analytical Batch: 271209

CONCENTRATION UNITS: μg/L

Analytical Method: OLCO 2.1

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|-----------|-------------------|--------|---|-------|-----|
| 127-18-4 | Tetrachloroethene | ND | U | 0.010 | 1.0 |
| 108-88-3 | Toluene | ND | U | 0.010 | 1.0 |
| 79-01-6 | Trichloroethene | ND | U | 0.010 | 1.0 |
| 75-01-4 | Vinyl chloride | ND | U | 0.010 | 1.0 |
| 1330-20-7 | Xylene tota | ND | U | 0.010 | 1.0 |

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

SKSW521009

Lab Name: GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204032408
Matrix: Water Lab Sample ID: 20403240806
Sample wt/vol: _____ Units: _____ Lab File ID: _____
Level: (low/med) _____ Date Collected: 03/22/04 Time: 1230
% Moisture: not dec. _____ Date Received: 03/24/04
GC Column: DB-624-30M ID: .25 (mm) Date Analyzed: 03/24/04 Time: 1801
Instrument ID: MSV2 Dilution Factor: 1 Analyst: RSP
Soil Extract Volume: _____ (μ L)
Soil Aliquot Volume: _____ (μ L)

Number TICs Found: 0

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | RT | EST. CONC. | Q |
|---------|------------------|----|------------|---|
| 1. | No tics detected | | | |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SKSW531009

Lab Name: GCAL

Contract

Lab Code: LA024

Case No.

SAS No

SDG No.: 204032408

Matrix: (soil/water) Water

Sample w/vol: 25 g/ml mL

Lab Sample ID: 20403240807

Level: (low/med)

Lab File ID: 2040324/T2585

% Moisture: not dec.

Date Collected: 03/22/04

Time: 1255

GC Column: DB-624-30M ID: 53 (mm)

Date Received: 03/24/04

Instrument ID: MSV2

Date Analyzed: 03/24/04

Time: 1826

Soil Extract Volume: (µL)

Dilution Factor: 1

Analyst: RJO

Soil Aliquot Volume: (µL)

Prep Batch:

Analytical Batch: 271209

CONCENTRATION UNITS ug/L

Analytical Method: OLC0 2.1

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|------------|---------------------------|--------|---|-------|-----|
| 71-55-6 | 1,1,1-Trichloroethane | 1.0 | U | 0.010 | 1.0 |
| 78-34-5 | 1,1,2,2-Tetrachloroethane | 1.0 | U | 0.010 | 1.0 |
| 78-00-5 | 1,1,2-Trichloroethane | 1.0 | U | 0.010 | 1.0 |
| 75-34-3 | 1,1-Dichloroethane | 1.0 | U | 0.010 | 1.0 |
| 75-35-4 | 1,1-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 106-93-4 | 1,2-Dibromoethane | 1.0 | U | 0.010 | 1.0 |
| 95-50-1 | 1,2-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 107-06-2 | 1,2-Dichloroethane | 1.0 | U | 0.010 | 1.0 |
| 540-59-0 | 1,2-Dichloroethene | 1.0 | U | 0.010 | 1.0 |
| 78-87-5 | 1,2-Dichloropropane | 1.0 | U | 0.010 | 1.0 |
| 541-73-1 | 1,3-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 106-46-7 | 1,4-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 78-93-3 | 2-Butanone | 5.0 | U | 0.010 | 5.0 |
| 501-78-6 | 2-Hexanone | 5.0 | U | 0.010 | 5.0 |
| 106-10-1 | 4-Methyl-2-pentanone | 5.0 | U | 0.010 | 5.0 |
| 67-64-1 | Acetone | 5.0 | U | 0.010 | 5.0 |
| 71-43-2 | Benzene | 1.0 | U | 0.010 | 1.0 |
| 75-27-4 | Bromodichloromethane | 1.0 | U | 0.010 | 1.0 |
| 75-25-2 | Bromoform | 1.0 | U | 0.010 | 1.0 |
| 74-83-9 | Bromomethane | 1.0 | U | 0.010 | 1.0 |
| 75-15-0 | Carbon disulfide | 1.0 | U | 0.010 | 1.0 |
| 56-23-5 | Carbon tetrachloride | 1.0 | U | 0.010 | 1.0 |
| 106-90-7 | Chlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 75-00-3 | Chloroethane | 1.0 | U | 0.010 | 1.0 |
| 67-66-3 | Chloroform | 1.0 | U | 0.010 | 1.0 |
| 74-87-3 | Chloromethane | 1.0 | U | 0.010 | 1.0 |
| 124-48-1 | Dibromochloromethane | 1.0 | U | 0.010 | 1.0 |
| 10061-01-5 | cis-1,3-Dichloropropene | 1.0 | U | 0.010 | 1.0 |
| 10061-02-6 | trans-1,3-Dichloropropene | 1.0 | U | 0.010 | 1.0 |
| 100-41-4 | Ethylbenzene | 1.0 | U | 0.010 | 1.0 |
| 75-09-2 | Methylene chloride | 2.0 | U | 0.010 | 2.0 |
| 100-42-5 | Styrene | 1.0 | U | 0.010 | 1.0 |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SKSW531009

Lab Name: GCAL Contract: _____

Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204032408

Matrix: (soil/water) Water

Sample wt/vol: 25 (g/ml) mL

Lab Sample ID: 20403240807

Level: (low/med) _____

Lab File ID: 2040324/T2585

% Moisture: not dec. _____

Date Collected: 03/22/04 Time: 1255

GC Column: DB-624-30M ID: .53 (mm)

Date Received: 03/24/04

Instrument ID: MSV2

Date Analyzed: 03/24/04 Time: 1826

Soil Extract Volume: _____ (µL)

Dilution Factor: 1 Analyst: RJO

Soil Aliquot Volume: _____ (µL)

Prep Batch: _____ Analytical Batch: 271209

CONCENTRATION UNITS: ug/L

Analytical Method: OLCO 2.1

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|-----------|-------------------|--------|---|-------|-----|
| 127-18-4 | Tetrachloroethene | 1.0 | U | 0.010 | 1.0 |
| 108-88-3 | Toluene | 1.0 | U | 0.010 | 1.0 |
| 79-01-6 | Trichloroethene | 1.0 | U | 0.010 | 1.0 |
| 75-01-4 | Vinyl chloride | 1.0 | U | 0.010 | 1.0 |
| 1330-20-7 | Xylene (total) | 1.0 | U | 0.010 | 1.0 |

VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

SKSW531009

Lab Name: GCAL Contract: _____
 Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204032408
 Matrix: Water Lab Sample ID: 20403240807
 Sample wt/vol: _____ Units: _____ Lab File ID: _____
 Level: (low/med) _____ Date Collected: 03/22/04 Time: 1255
 % Moisture: not dec. _____ Date Received: 03/24/04
 GC Column: DB-624-30M ID: 25 (mm) Date Analyzed: 03/24/04 Time: 1826
 Instrument ID: MSV2 Dilution Factor: 1 Analyst: RSP
 Soil Extract Volume: _____ (µL)
 Soil Aliquot Volume: _____ (µL)

Number TICs Found: 0

CONCENTRATION UNITS

| CAS NO. | COMPOUND | RT | EST. CONC. | Q |
|---------|------------------|----|------------|---|
| 1. | No tics detected | | | |

SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

Lab Name: GCAL Sample ID: SKSWEB1009
 Lab Code: LA024 Case No.: _____ Contract: _____
 SAS No.: _____ SDG No.: 204032408 Lab File ID: 2040407P/S5094
 Matrix: Water Lab Sample ID: 20403240801
 Sample wt/vol: 1000 Units: mL Date Collected: 03/22/04 Time: 1345
 Level: (low/med) _____ Date Received: 03/24/04
 % Moisture: _____ decanted: (Y/N) _____ Date Extracted: 3/25/04
 GC Column: DB-5MS-30M ID: .25 (mm) Date Analyzed: 04/07/04 Time: 1838
 Concentrated Extract Volume: 1000 (µL) Dilution Factor: 1 Analyst: RLW
 Injection Volume: 1.0 (µL) Prep Method: _____
 GPC Cleanup: (Y/N) N pH: _____ Analytical Method: OLMO 4.2
 Instrument ID: MSSV2

CONCENTRATION UNITS: ug/LPrep Batch: 271238 Analytical Batch: 271976

| CAS NO. | COMPOUND | RESULT | Q | MDL | PQL |
|-----------|-----------------------------|--------|---|-------|------|
| 95-95-4 | 2,4,5-Trichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 88-06-2 | 2,4,6-Trichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 120-83-2 | 2,4-Dichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 51-28-5 | 2,4-Dinitrophenol | 25.0 | U | 0.010 | 25.0 |
| 121-14-2 | 2,4-Dinitrotoluene | 10.0 | U | 0.010 | 10.0 |
| 606-20-2 | 2,6-Dinitrotoluene | 10.0 | U | 0.010 | 10.0 |
| 91-58-7 | 2-Chloronaphthalene | 10.0 | U | 0.010 | 10.0 |
| 95-57-8 | 2-Chlorophenol | 10.0 | U | 0.010 | 10.0 |
| 91-57-6 | 2-Methylnaphthalene | 10.0 | U | 0.010 | 10.0 |
| 88-74-4 | 2-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 88-75-5 | 2-Nitrophenol | 10.0 | U | 0.010 | 10.0 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 10.0 | U | 0.010 | 10.0 |
| 99-09-2 | 3-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 534-52-1 | 2-Methyl-4,6-dinitrophenol | 25.0 | U | 0.010 | 25.0 |
| 59-50-7 | 4-Chloro-3-methylphenol | 10.0 | U | 0.010 | 10.0 |
| 106-47-8 | 4-Chloroaniline | 10.0 | U | 0.010 | 10.0 |
| 7005-72-3 | 4-Chlorophenyl-phenylether | 10.0 | U | 0.010 | 10.0 |
| 106-44-5 | 4-Methylphenol (p-Cresol) | 10.0 | U | 0.010 | 10.0 |
| 83-32-9 | Acenaphthene | 10.0 | U | 0.010 | 10.0 |
| 208-96-8 | Acenaphthylene | 10.0 | U | 0.010 | 10.0 |
| 120-12-7 | Anthracene | 10.0 | U | 0.010 | 10.0 |
| 56-55-3 | Benzo(a)anthracene | 10.0 | U | 0.010 | 10.0 |
| 50-32-8 | Benzo(a)pyrene | 10.0 | U | 0.010 | 10.0 |
| 205-99-2 | Benzo(b)fluoranthene | 10.0 | U | 0.010 | 10.0 |
| 191-24-2 | Benzo(g,h,i)perylene | 10.0 | U | 0.010 | 10.0 |
| 207-08-9 | Benzo(k)fluoranthene | 10.0 | U | 0.010 | 10.0 |
| 111-91-1 | Bis(2-Chloroethoxy)methane | 10.0 | U | 0.010 | 10.0 |
| 111-44-4 | Bis(2-Chloroethyl)ether | 10.0 | U | 0.010 | 10.0 |
| 108-60-1 | bis(2-Chloroisopropyl)ether | 10.0 | U | 0.010 | 10.0 |
| 117-81-7 | bis(2-ethylhexyl)phthalate | 10.0 | U | 0.010 | 10.0 |

SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

Lab Name: GCAL Sample ID: SKSWEB1009
 Lab Code: LA024 Case No.: _____ Contract: _____
 SAS No.: _____ SOG No.: 204032408 Lab File ID: 2040407P/S5094
 Matrix: Water Lab Sample ID: 20403240801
 Sample wt/vol: 1000 Units: mL Date Collected: 03/22/04 Time: 1345
 Level: (low/med) _____ Date Received: 03/24/04
 % Moisture: _____ decanted: (Y/N) _____ Date Extracted: 3/25/04
 GC Column: DB-5MS-30M ID: 25 (mm) Date Analyzed: 04/07/04 Time: 1838
 Concentrated Extract Volume: 1000 (µL) Dilution Factor: 1 Analyst: RLW
 Injection Volume: 10 (µL) Prep Method: _____
 GPC Cleanup: (Y/N) N pH: _____ Analytical Method: OLMD 4.2
 Instrument ID: MSSV2

CONCENTRATION UNITS: ug/L

Prep Batch: 271238 Analytical Batch: 271976

| CAS NO. | COMPOUND | RESULT | Q | MDL | PQL |
|----------|---------------------------|--------|---|-------|------|
| 101-55-3 | 4-Bromophenyl-phenylether | 10.0 | U | 0.010 | 10.0 |
| 85-68-7 | Butylbenzylchloralate | 10.0 | U | 0.010 | 10.0 |
| 86-74-8 | Carbazole | 10.0 | U | 0.010 | 10.0 |
| 218-01-9 | Chrysene | 10.0 | U | 0.010 | 10.0 |
| 84-74-2 | Di-n-butylphthalate | 10.0 | U | 0.010 | 10.0 |
| 117-84-0 | Di-n-octylphthalate | 10.0 | U | 0.010 | 10.0 |
| 53-70-3 | Dibenz(a,h)anthracene | 10.0 | U | 0.010 | 10.0 |
| 132-64-9 | Dibenzofuran | 10.0 | U | 0.010 | 10.0 |
| 84-86-2 | Diethylphthalate | 10.0 | U | 0.010 | 10.0 |
| 131-11-3 | Dimethyl-phthalate | 10.0 | U | 0.010 | 10.0 |
| 105-67-9 | 2,4-Dimethylphenol | 10.0 | U | 0.010 | 10.0 |
| 208-44-0 | Fluoranthene | 10.0 | U | 0.010 | 10.0 |
| 86-73-7 | Fluorene | 10.0 | U | 0.010 | 10.0 |
| 118-74-1 | Hexachlorobenzene | 10.0 | U | 0.010 | 10.0 |
| 87-68-3 | Hexachlorobutadiene | 10.0 | U | 0.010 | 10.0 |
| 77-47-4 | Hexachlorocyclopentadiene | 10.0 | U | 0.010 | 10.0 |
| 67-72-1 | Hexachloroethane | 10.0 | U | 0.010 | 10.0 |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | 10.0 | U | 0.010 | 10.0 |
| 78-59-1 | Isophorone | 10.0 | U | 0.010 | 10.0 |
| 91-20-3 | Naphthalene | 10.0 | U | 0.010 | 10.0 |
| 100-01-6 | 4-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 98-95-3 | Nitrobenzene | 10.0 | U | 0.010 | 10.0 |
| 100-02-7 | 4-Nitrophenol | 25.0 | U | 0.010 | 25.0 |
| 87-86-5 | Pentachlorophenol | 25.0 | U | 0.010 | 25.0 |
| 85-01-8 | Phenanthrene | 10.0 | U | 0.010 | 10.0 |
| 108-95-2 | Phenol | 10.0 | U | 0.010 | 10.0 |
| 129-00-0 | Pyrene | 10.0 | U | 0.010 | 10.0 |
| 621-64-7 | N-Nitroso-dim-propylamine | 10.0 | U | 0.010 | 10.0 |
| 86-30-6 | N-Nitrosodiphenylamine | 10.0 | U | 0.010 | 10.0 |
| 95-48-7 | o-Cresol | 10.0 | U | 0.010 | 10.0 |

SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

| | |
|---|---|
| Lab Name: <u>GCAL</u> | Sample ID: <u>SKSWEB1009</u> |
| Lab Code: <u>LA024</u> <u>2</u> Case No.: _____ | Contract: _____ |
| SAS No.: _____ SDG No.: <u>204032408</u> | Lab File ID: <u>S5094</u> |
| Matrix: <u>Water</u> | Lab Sample ID: <u>20403240801</u> |
| Sample wt/vol: _____ Units: _____ | Date Collected: <u>03/22/04</u> Time: <u>1345</u> |
| Level: (low/med) _____ | Date Received: <u>03/24/04</u> |
| % Moisture: not dec. _____ | Date Extracted: <u>3/25/04</u> |
| GC Column: _____ ID: _____ (mm) | Date Analyzed: <u>04/07/04</u> Time: <u>1838</u> |
| Concentrated Extract Volume: <u>1000</u> (µL) | Dilution Factor: <u>1</u> Analyst: <u>RLW</u> |
| Injection Volume: <u>1.0</u> (µL) | Prep Method: _____ |
| GPC Cleanup: (Y/N) <u>N</u> pH: _____ | Analytical Method: <u>SW-846 8270C</u> |
| | Instrument ID: <u>MSSV2</u> |

Number TICs Found : 0

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | RT | EST. CONC. | Q |
|---------|------------------|----|------------|---|
| 1. | No tics detected | | | |

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SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

Lab Name: GCAL
 Lab Code: LA024 Case No.: _____
 SAS No.: _____ SDG No.: 204032408
 Matrix: Water
 Sample wt/vol: 1000 Units mL
 Level: (low/med) _____
 % Moisture: _____ decahed: (Y/N) _____
 GC Column: DB-5MS-30M ID: 25 (mm)
 Concentrated Extract Volume: 1000 (µL)
 Injection Volume: 1.0 (µL)
 GPC Cleanup: (Y/N) N pH: _____

Sample ID: SKSW511009
 Contract: _____
 Lab File ID: 2040407P/S5095
 Lab Sample ID: 20403240802
 Date Collected: 03/22/04 Time: 1040
 Date Received: 03/24/04
 Date Extracted: 3/25/04
 Date Analyzed: 04/07/04 Time: 1905
 Dilution Factor: 1 Analyst: RLW
 Prep Method: _____
 Analytical Method: OLMO 4.2
 Instrument ID: MSSV2

CONCENTRATION UNITS: ug/L

Prep Batch: 271238 Analytical Batch: 271976

| CAS NO. | COMPOUND | RESULT | Q | MDL | PQL |
|-----------|-----------------------------|--------|---|-------|------|
| 95-95-4 | 2,4,5-Trichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 98-06-2 | 2,4,6-Trichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 120-83-2 | 2,4-Dichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 51-28-5 | 2,4-Dinitrophenol | 25.0 | U | 0.010 | 25.0 |
| 121-14-2 | 2,4-Dinitrotoluene | 10.0 | U | 0.010 | 10.0 |
| 608-20-2 | 2,6-Dinitrotoluene | 10.0 | U | 0.010 | 10.0 |
| 91-58-7 | 2-Chloronaphthalene | 10.0 | U | 0.010 | 10.0 |
| 95-57-8 | 2-Chlorophenol | 10.0 | U | 0.010 | 10.0 |
| 91-57-6 | 2-Methylnaphthalene | 10.0 | U | 0.010 | 10.0 |
| 88-74-4 | 2-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 88-75-5 | 2-Nitrophenol | 10.0 | U | 0.010 | 10.0 |
| 91-84-1 | 3,3'-Dichlorobenzidine | 10.0 | U | 0.010 | 10.0 |
| 99-09-2 | 3-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 534-52-1 | 2-Methyl-4,5-dinitrophenol | 25.0 | U | 0.010 | 25.0 |
| 59-50-7 | 4-Chloro-3-methylphenol | 10.0 | U | 0.010 | 10.0 |
| 108-47-8 | 4-Chloroaniline | 10.0 | U | 0.010 | 10.0 |
| 7005-72-3 | 4-Chlorophenylphenylether | 10.0 | U | 0.010 | 10.0 |
| 108-44-5 | 4-Methylphenol (p-Cresol) | 10.0 | U | 0.010 | 10.0 |
| 63-32-9 | Acenaphthene | 10.0 | U | 0.010 | 10.0 |
| 208-96-8 | Acenaphthylene | 10.0 | U | 0.010 | 10.0 |
| 120-12-7 | Anthracene | 10.0 | U | 0.010 | 10.0 |
| 58-55-3 | Benzofluoranthene | 10.0 | U | 0.010 | 10.0 |
| 50-32-6 | Benzofluoranthene | 10.0 | U | 0.010 | 10.0 |
| 205-99-2 | Benzofluoranthene | 10.0 | U | 0.010 | 10.0 |
| 191-24-2 | Benzofluoranthene | 10.0 | U | 0.010 | 10.0 |
| 207-08-9 | Benzofluoranthene | 10.0 | U | 0.010 | 10.0 |
| 111-91-1 | Bis(2-Chloroethoxy)methane | 10.0 | U | 0.010 | 10.0 |
| 111-44-4 | Bis(2-Chloroethyl)ether | 10.0 | U | 0.010 | 10.0 |
| 108-60-1 | bis(2-Chloroisopropyl)ether | 10.0 | U | 0.010 | 10.0 |
| 117-81-7 | bis(2-ethylhexyl)phthalate | 10.0 | U | 0.010 | 10.0 |

SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

Lab Name: GCAL Sample ID: SKSW511009
 Lab Code: LA024 Case No.: _____ Contract: _____
 SAS No.: _____ SDG No.: 204032408 Lab File ID: 2040407P/S5095
 Matrix: Water Lab Sample ID: 20403240802
 Sample wt/vol: 1000 Units: mL Date Collected: 03/22/04 Time: 1040
 Level: (low/med) _____ Date Received: 03/24/04
 % Moisture: _____ decanted: (Y/N) _____ Date Extracted: 3/25/04
 GC Column: DB-5MS-30M ID: .25 (mm) Date Analyzed: 04/07/04 Time: 1905
 Concentrated Extract Volume: 1000 (µL) Dilution Factor: 1 Analyst: RLW
 Injection Volume: 1.0 (µL) Prep Method: _____
 GPC Cleanup: (Y/N) N pH: _____ Analytical Method: OLMO 4.2
 Instrument ID: MSSV2
 CONCENTRATION UNITS: ug/L Prep Batch: 271238 Analytical Batch: 271976

| CAS NO. | COMPOUND | RESULT | Q | MDL | PQL |
|----------|----------------------------|--------|---|-------|------|
| 101-55-3 | 4-Bromophenyl-phenylether | 10.0 | U | 0.010 | 10.0 |
| 85-68-7 | Butylbenzylphthalate | 10.0 | U | 0.010 | 10.0 |
| 86-74-8 | Carbazole | 10.0 | U | 0.010 | 10.0 |
| 218-01-9 | Chrysene | 10.0 | U | 0.010 | 10.0 |
| 84-74-2 | Di-n-butylphthalate | 10.0 | U | 0.010 | 10.0 |
| 117-84-0 | Di-n-octylphthalate | 10.0 | U | 0.010 | 10.0 |
| 53-70-3 | Dibenz(a,h)anthracene | 10.0 | U | 0.010 | 10.0 |
| 132-64-9 | Dibenzofuran | 10.0 | U | 0.010 | 10.0 |
| 84-66-2 | Diethylphthalate | 10.0 | U | 0.010 | 10.0 |
| 131-11-3 | Dimethyl-phthalate | 10.0 | U | 0.010 | 10.0 |
| 105-67-9 | 2,4-Dimethylphenol | 10.0 | U | 0.010 | 10.0 |
| 206-44-0 | Fluoranthene | 0.849 | J | 0.010 | 10.0 |
| 86-73-7 | Fluorene | 10.0 | U | 0.010 | 10.0 |
| 118-74-1 | Hexachlorobenzene | 10.0 | U | 0.010 | 10.0 |
| 87-68-3 | Hexachlorobutadiene | 10.0 | U | 0.010 | 10.0 |
| 77-47-4 | Hexachlorocyclopentadiene | 10.0 | U | 0.010 | 10.0 |
| 67-72-1 | Hexachloroethane | 10.0 | U | 0.010 | 10.0 |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | 10.0 | U | 0.010 | 10.0 |
| 78-59-1 | Isophorone | 10.0 | U | 0.010 | 10.0 |
| 91-20-3 | Naphthalene | 10.0 | U | 0.010 | 10.0 |
| 100-01-6 | 4-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 98-95-3 | Nitrobenzene | 10.0 | U | 0.010 | 10.0 |
| 100-02-7 | 4-Nitrophenol | 25.0 | U | 0.010 | 25.0 |
| 87-86-5 | Pentachlorophenol | 25.0 | U | 0.010 | 25.0 |
| 85-01-8 | Phenanthrene | 0.794 | J | 0.010 | 10.0 |
| 108-95-2 | Phenol | 10.0 | U | 0.010 | 10.0 |
| 129-00-0 | Pyrene | 0.679 | J | 0.010 | 10.0 |
| 621-64-7 | N-Nitroso-di-n-propylamine | 10.0 | U | 0.010 | 10.0 |
| 86-30-6 | N-Nitrosodiphenylamine | 10.0 | U | 0.010 | 10.0 |
| 95-48-7 | o-Cresol | 10.0 | U | 0.010 | 10.0 |

43
 43

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SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

| | |
|---|---|
| Lab Name: <u>GCAL</u> | Sample ID: <u>SKSW511009</u> |
| Lab Code: <u>LA024</u> <u>2</u> Case No.: _____ | Contract: _____ |
| SAS No.: _____ SDG No.: <u>204032408</u> | Lab File ID: <u>S5095</u> |
| Matrix: <u>Water</u> | Lab Sample ID: <u>20403240802</u> |
| Sample wt/vol: _____ Units: _____ | Date Collected: <u>03/22/04</u> Time: <u>1040</u> |
| Level: (low/med) _____ | Date Received: <u>03/24/04</u> |
| % Moisture: not dec. _____ | Date Extracted: <u>3/25/04</u> |
| GC Column: _____ ID: _____ (mm) | Date Analyzed: <u>04/07/04</u> Time: <u>1905</u> |
| Concentrated Extract Volume: <u>1000</u> (µL) | Dilution Factor: <u>1</u> Analyst: <u>RLW</u> |
| Injection Volume: <u>10</u> (µL) | Prep Method: _____ |
| GPC Cleanup: (Y/N) <u>N</u> pH: _____ | Analytical Method: <u>SW-846 8270C</u> |
| | Instrument ID: <u>MSSV2</u> |

Number TICs Found: 0

CONCENTRATION UNITS

| CAS NO. | COMPOUND | RT | EST. CONC. | Q |
|---------|------------------|----|------------|---|
| 1. | No TICs detected | | | |

5/26/04
RLW

SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

Lab Name: GCAL Sample ID: SKSW521009
 Lab Code: LA024 Case No.: _____ Contract: _____
 SAS No.: _____ SDG No.: 204032408 Lab File ID: 2040407P/S5098
 Matrix: Water Lab Sample ID: 20403240806
 Sample wt/vol: 1000 Units: mL Date Collected: 03/22/04 Time: 1230
 Level: (low/med) _____ Date Received: 03/24/04
 % Moisture: _____ decanted: (Y/N) _____ Date Extracted: 3125104
 GC Column: DB-5MS-30M ID: .25 (mm) Date Analyzed: 04/07/04 Time: 2025
 Concentrated Extract Volume: 1000 (µL) Dilution Factor: 1 Analyst: RLW
 Injection Volume: 1.0 (µL) Prep Method: _____
 GPC Cleanup: (Y/N) N pH: _____ Analytical Method: OLMO 4.2
 Instrument ID: MSSV2
 CONCENTRATION UNITS: ug/L Prep Batch: 271238 Analytical Batch: 271976

| CAS NO. | COMPOUND | RESULT | Q | MDL | PQL |
|-----------|-----------------------------|--------|---|-------|------|
| 95-95-4 | 2,4,5-Trichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 88-06-2 | 2,4,6-Trichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 120-83-2 | 2,4-Dichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 51-28-5 | 2,4-Dinitrophenol | 25.0 | U | 0.010 | 25.0 |
| 121-14-2 | 2,4-Dinitrotoluene | 10.0 | U | 0.010 | 10.0 |
| 606-20-2 | 2,6-Dinitrotoluene | 10.0 | U | 0.010 | 10.0 |
| 91-58-7 | 2-Chloronaphthalene | 10.0 | U | 0.010 | 10.0 |
| 95-57-8 | 2-Chlorophenol | 10.0 | U | 0.010 | 10.0 |
| 91-57-6 | 2-Methylnaphthalene | 10.0 | U | 0.010 | 10.0 |
| 88-74-4 | 2-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 88-75-5 | 2-Nitrophenol | 10.0 | U | 0.010 | 10.0 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 10.0 | U | 0.010 | 10.0 |
| 99-09-2 | 3-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 534-52-1 | 2-Methyl-4,6-dinitrophenol | 25.0 | U | 0.010 | 25.0 |
| 59-50-7 | 4-Chloro-3-methylphenol | 10.0 | U | 0.010 | 10.0 |
| 106-47-8 | 4-Chloroaniline | 10.0 | U | 0.010 | 10.0 |
| 7005-72-3 | 4-Chlorophenyl-phenylether | 10.0 | U | 0.010 | 10.0 |
| 106-44-5 | 4-Methylphenol (p-Cresol) | 10.0 | U | 0.010 | 10.0 |
| 83-32-9 | Acenaphthene | 10.0 | U | 0.010 | 10.0 |
| 208-96-8 | Acenaphthylene | 10.0 | U | 0.010 | 10.0 |
| 120-12-7 | Anthracene | 10.0 | U | 0.010 | 10.0 |
| 56-55-3 | Benzo(a)anthracene | 10.0 | U | 0.010 | 10.0 |
| 50-32-8 | Benzo(a)pyrene | 10.0 | U | 0.010 | 10.0 |
| 205-99-2 | Benzo(b)fluoranthene | 10.0 | U | 0.010 | 10.0 |
| 191-24-2 | Benzo(g,h,i)perylene | 10.0 | U | 0.010 | 10.0 |
| 207-08-9 | Benzo(k)fluoranthene | 10.0 | U | 0.010 | 10.0 |
| 111-91-1 | Bis(2-Chloroethoxy)methane | 10.0 | U | 0.010 | 10.0 |
| 111-44-4 | Bis(2-Chloroethyl)ether | 10.0 | U | 0.010 | 10.0 |
| 108-60-1 | bis(2-Chloroisopropyl)ether | 10.0 | U | 0.010 | 10.0 |
| 117-81-7 | bis(2-ethylhexyl)phthalate | 10.0 | U | 0.010 | 10.0 |

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SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

Lab Name: GCAL Sample ID: SKSW521009
 Lab Code: LA024 Case No: _____ Contract: _____
 SAS No.: _____ SDG No.: 204032408 Lab File ID: 2040407P/S5098
 Matrix: Water Lab Sample ID: 20403240806
 Sample wt/vol: 1000 units mL Date Collected: 03/22/04 Time: 1230
 Level: (low/med) _____ Date Received: 03/24/04
 % Moisture: _____ decanted (Y/N) _____ Date Extracted: 3/25/04
 GC Column: DB-5MS-30M ID: 25 (mm) Date Analyzed: 04/07/04 Time: 2025
 Concentrated Extract Volume: 1000 (µL) Dilution Factor: 1 Analyst: RLW
 Injection Volume: 1.0 (µL) Prep Method: _____
 GPC Cleanup: (Y/N) N pH: _____ Analytical Method: OLMO 4.2

CONCENTRATION UNITS µg/L

| CAS NO. | COMPOUND | RESULT | Q | MDL | PQL |
|----------|----------------------------|--------|---|-------|------|
| 101-55-3 | 4-Bromophenyl-phenylether | 10.0 | U | 0.010 | 10.0 |
| 85-68-7 | Butylbenzylphthalate | 10.0 | U | 0.010 | 10.0 |
| 86-74-8 | Carbazole | 10.0 | U | 0.010 | 10.0 |
| 218-01-9 | Chrysene | 10.0 | U | 0.010 | 10.0 |
| 84-74-2 | Di-n-butylphthalate | 10.0 | U | 0.010 | 10.0 |
| 117-84-0 | Di-n-octylphthalate | 10.0 | U | 0.010 | 10.0 |
| 53-70-3 | Dibenz(a,h)anthracene | 10.0 | U | 0.010 | 10.0 |
| 132-64-9 | Dibenzofuran | 10.0 | U | 0.010 | 10.0 |
| 84-66-2 | Diethylphthalate | 10.0 | U | 0.010 | 10.0 |
| 131-11-3 | Dimethyl-phthalate | 10.0 | U | 0.010 | 10.0 |
| 105-67-8 | 2,4-Dimethylphenol | 10.0 | U | 0.010 | 10.0 |
| 208-44-0 | Fluoranthene | 10.0 | U | 0.010 | 10.0 |
| 86-73-7 | Fluorene | 10.0 | U | 0.010 | 10.0 |
| 118-74-1 | Hexachlorobenzene | 10.0 | U | 0.010 | 10.0 |
| 87-68-3 | Hexachlorobutadiene | 10.0 | U | 0.010 | 10.0 |
| 77-47-4 | Hexachlorocyclopentadiene | 10.0 | U | 0.010 | 10.0 |
| 67-72-1 | Hexachloroethane | 10.0 | U | 0.010 | 10.0 |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | 10.0 | U | 0.010 | 10.0 |
| 78-58-1 | Isophorone | 10.0 | U | 0.010 | 10.0 |
| 91-20-3 | Naphthalene | 10.0 | U | 0.010 | 10.0 |
| 100-01-6 | 4-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 98-95-3 | Nitrobenzene | 10.0 | U | 0.010 | 10.0 |
| 100-02-7 | 4-Nitrophenol | 25.0 | U | 0.010 | 25.0 |
| 87-86-5 | Pentachlorobenzene | 25.0 | U | 0.010 | 25.0 |
| 85-01-8 | Phenanthrene | 10.0 | U | 0.010 | 10.0 |
| 108-95-2 | Phenol | 10.0 | U | 0.010 | 10.0 |
| 129-00-0 | Pyrene | 10.0 | U | 0.010 | 10.0 |
| 621-84-7 | N-Nitroso-di-n-propylamine | 10.0 | U | 0.010 | 10.0 |
| 86-30-6 | N-Nitrosodiphenylamine | 10.0 | U | 0.010 | 10.0 |
| 95-48-7 | o-Cresol | 10.0 | U | 0.010 | 10.0 |

SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

Lab Name: GCAL Sample ID: SKSW521009
Lab Code: LA024 2 Case No.: _____ Contract: _____
SAS No.: _____ SDG No.: 204032408 Lab File ID: S5098
Matrix: Water Lab Sample ID: 20403240806
Sample wt/vol: _____ Units: _____ Date Collected: 03/22/04 Time: 1230
Level: (low/med) _____ Date Received: 03/24/04
% Moisture: not dec. _____ Date Extracted: 3/25/04
GC Column: _____ ID: _____ (mm) Date Analyzed: 04/07/04 Time: 2025
Concentrated Extract Volume: 1000 (µL) Dilution Factor: 1 Analyst: RLW
Injection Volume: 1.0 (µL) Prep Method: _____
GPC Cleanup: (Y/N) N pH: _____ Analytical Method: SW-846 8270C
Instrument ID: MSSV2

Number TICs Found : 0

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | RT | EST. CONC. | Q |
|---------|------------------|----|------------|---|
| 1. | No tics detected | | | |

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SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

Lab Name: GCAL Sample ID: SKSW531009
 Lab Code: LA024 Case No: _____ Contract: _____
 SAS No.: _____ SDG No.: 204032408 Lab File ID: 2040407P/S5099
 Matrix: Water Lab Sample ID: 20403240807
 Sample w/voit: 1000 units mL Date Collected: 03/22/04 Time: 1255
 Level: (low/med) _____ Date Received: 03/24/04
 % Moisture: _____ decanted: (Y/N) _____ Date Extracted: 3/25/04
 GC Column: DB-5MS-30M ID: 25 (mm) Date Analyzed: 04/07/04 Time: 2053
 Concentrated Extract Volume: 1000 (µL) Dilution Factor: 1 Analyst: RLW
 Injection Volume: 1.0 (µL) Prep Method: _____
 GPC Cleanup: (Y/N) N pH: _____ Analytical Method: OLMO 4.2

CONCENTRATION UNITS µg/L

Instrument ID: MSSV2
 Prep Batch: 271238 Analytical Batch: 271976

| CAS NO. | COMPOUND | RESULT | Q | MDL | PQL |
|-----------|-----------------------------|--------|---|-------|------|
| 95-95-4 | 2,4,5-Trichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 88-06-2 | 2,4,6-Trichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 120-83-2 | 2,4-Dichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 51-28-5 | 2,4-Dinitrophenol | 25.0 | U | 0.010 | 25.0 |
| 121-14-2 | 2,4-Dinitrotoluene | 10.0 | U | 0.010 | 10.0 |
| 806-20-2 | 2,6-Dinitrotoluene | 10.0 | U | 0.010 | 10.0 |
| 91-58-7 | 2-Chloronaphthalene | 10.0 | U | 0.010 | 10.0 |
| 95-57-8 | 2-Chlorophenol | 10.0 | U | 0.010 | 10.0 |
| 91-57-6 | 2-Methylnaphthalene | 10.0 | U | 0.010 | 10.0 |
| 88-74-4 | 2-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 88-75-5 | 2-Nitrophenol | 10.0 | U | 0.010 | 10.0 |
| 91-04-1 | 3,3'-Dichlorobenzidine | 10.0 | U | 0.010 | 10.0 |
| 99-08-2 | 3-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 534-52-1 | 2-Methyl-4,6-dinitrophenol | 25.0 | U | 0.010 | 25.0 |
| 59-50-7 | 4-Chloro-3-methylphenol | 10.0 | U | 0.010 | 10.0 |
| 105-47-8 | 4-Chloroaniline | 10.0 | U | 0.010 | 10.0 |
| 7005-72-3 | 4-Chlorophenyl-phenylether | 10.0 | U | 0.010 | 10.0 |
| 108-44-5 | 4-Methylphenol (p-Cresol) | 10.0 | U | 0.010 | 10.0 |
| 83-32-9 | Acenaphthene | 10.0 | U | 0.010 | 10.0 |
| 208-96-8 | Acenaphthylene | 10.0 | U | 0.010 | 10.0 |
| 120-12-7 | Anthracene | 10.0 | U | 0.010 | 10.0 |
| 56-55-3 | Benzo(a)anthracene | 10.0 | U | 0.010 | 10.0 |
| 50-32-8 | Benzo(a)pyrene | 10.0 | U | 0.010 | 10.0 |
| 205-99-2 | Benzo(b)fluoranthene | 10.0 | U | 0.010 | 10.0 |
| 191-24-2 | Benzo(g,h,i)perylene | 10.0 | U | 0.010 | 10.0 |
| 207-08-9 | Benzo(k)fluoranthene | 10.0 | U | 0.010 | 10.0 |
| 111-91-1 | Bis(2-Chloroethoxy)methane | 10.0 | U | 0.010 | 10.0 |
| 111-44-4 | Bis(2-Chloroethyl)ether | 10.0 | U | 0.010 | 10.0 |
| 108-60-1 | bis(2-Chloroisopropyl)ether | 10.0 | U | 0.010 | 10.0 |
| 117-81-7 | bis(2-ethylhexyl) sebacate | 10.0 | U | 0.010 | 10.0 |

3/24/04
RLW

SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

Lab Name: GCAL Sample ID: SKSW531009
 Lab Code: LA024 Case No.: _____ Contract: _____
 SAS No.: _____ SDG No.: 204032408 Lab File ID: 2040407P/S5099
 Matrix: Water Lab Sample ID: 20403240807
 Sample wt/vol: 1000 Units: mL Date Collected: 03/22/04 Time: 1255
 Level: (low/med) _____ Date Received: 03/24/04
 % Moisture: _____ decanted: (Y/N) _____ Date Extracted: 3/25/04
 GC Column: DB-5MS-30M ID: .25 (mm) Date Analyzed: 04/07/04 Time: 2053
 Concentrated Extract Volume: 1000 (µL) Dilution Factor: 1 Analyst: RLW
 Injection Volume: 1.0 (µL) Prep Method: _____
 GPC Cleanup: (Y/N) N pH: _____ Analytical Method: OLMO 4.2
 Instrument ID: MSSV2

CONCENTRATION UNITS: ug/LPrep Batch: 271238 Analytical Batch: 271976

| CAS NO. | COMPOUND | RESULT | Q | MDL | PQL |
|----------|----------------------------|--------|---|-------|------|
| 101-55-3 | 4-Bromophenyl-phenylether | 10.0 | U | 0.010 | 10.0 |
| 85-68-7 | Butylbenzylphthalate | 10.0 | U | 0.010 | 10.0 |
| 86-74-8 | Carbazole | 10.0 | U | 0.010 | 10.0 |
| 218-01-9 | Chrysene | 10.0 | U | 0.010 | 10.0 |
| 84-74-2 | Di-n-butylphthalate | 10.0 | U | 0.010 | 10.0 |
| 117-84-0 | Di-n-octylphthalate | 10.0 | U | 0.010 | 10.0 |
| 53-70-3 | Dibenz(a,h)anthracene | 10.0 | U | 0.010 | 10.0 |
| 132-64-9 | Dibenzofuran | 10.0 | U | 0.010 | 10.0 |
| 84-66-2 | Diethylphthalate | 10.0 | U | 0.010 | 10.0 |
| 131-11-3 | Dimethyl-phthalate | 10.0 | U | 0.010 | 10.0 |
| 105-67-9 | 2,4-Dimethylphenol | 10.0 | U | 0.010 | 10.0 |
| 206-44-0 | Fluoranthene | 10.0 | U | 0.010 | 10.0 |
| 86-73-7 | Fluorene | 10.0 | U | 0.010 | 10.0 |
| 113-74-1 | Hexachlorobenzene | 10.0 | U | 0.010 | 10.0 |
| 87-68-3 | Hexachlorobutadiene | 10.0 | U | 0.010 | 10.0 |
| 77-47-4 | Hexachlorocyclopentadiene | 10.0 | U | 0.010 | 10.0 |
| 67-72-1 | Hexachloroethane | 10.0 | U | 0.010 | 10.0 |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | 10.0 | U | 0.010 | 10.0 |
| 78-59-1 | Isophorone | 10.0 | U | 0.010 | 10.0 |
| 91-20-3 | Naphthalene | 10.0 | U | 0.010 | 10.0 |
| 100-01-6 | 4-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 98-95-3 | Nitrobenzene | 10.0 | U | 0.010 | 10.0 |
| 100-02-7 | 4-Nitrophenol | 25.0 | U | 0.010 | 25.0 |
| 87-86-5 | Pentachlorophenol | 25.0 | U | 0.010 | 25.0 |
| 85-01-8 | Phenanthrene | 10.0 | U | 0.010 | 10.0 |
| 108-95-2 | Phenol | 10.0 | U | 0.010 | 10.0 |
| 129-00-0 | Pyrene | 10.0 | U | 0.010 | 10.0 |
| 621-64-7 | N-Nitroso-di-n-propylamine | 10.0 | U | 0.010 | 10.0 |
| 86-30-6 | N-Nitrosodiphenylamine | 10.0 | U | 0.010 | 10.0 |
| 95-48-7 | o-Cresol | 10.0 | U | 0.010 | 10.0 |

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SEMI-VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

| | |
|--|---|
| Lab Name: <u>GCAL</u> | Sample ID: <u>SKSW531009</u> |
| Lab Code: <u>LA024</u> <u>2</u> Case No.: | Contract: |
| SAS No.: | Lab File ID: <u>S5099</u> |
| SDG No. <u>204032408</u> | Lab Sample ID: <u>20403240807</u> |
| Matrix: <u>Water</u> | Date Collected: <u>03/22/04</u> Time: <u>1255</u> |
| Sample wt/vol: _____ Units: _____ | Date Received: <u>03/24/04</u> |
| Level: (low/med) _____ | Date Extracted: <u>3/25/04</u> |
| % Moisture: not dec. _____ | Date Analyzed: <u>04/07/04</u> Time: <u>2053</u> |
| GC Column: _____ ID: _____ (mm) | Dilution Factor: <u>1</u> Analyst: <u>RLW</u> |
| Concentrated Extract Volume <u>1000</u> (μL) | Prep Method: _____ |
| Injection Volume: _____ (μL) | Analytical Method: <u>SW-846 8270C</u> |
| GPC Cleanup: (Y/N) <u>N</u> pH: _____ | Instrument ID: <u>MSSV2</u> |

Number TICs Found: 0

CONCENTRATION UNITS

| CAS NO. | COMPOUND | RT | EST. CONC. | Q |
|---------|-----------------|----|------------|---|
| 1. | No bcs detected | | | |

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1D
ORGANICS ANALYSIS DATA SHEET

Lab Name: GCAL Sample ID: SKSWEB1009
 Lab Code: LA024 Case No.: _____ Contract: _____
 Matrix: Water SAS No.: _____ SDG No.: 204032408
 Sample wt/vol: 1000 Units: mL Lab Sample ID: 20403240801
 Level: (low/med) _____ Date Collected: 03/22/04 Time: 1345
 % Moisture: _____ decanted: (Y/N) _____ Date Received: 03/24/04
 GC Column: DB-608-30M ID: .53 (mm) Date Extracted: 3/26/04
 Concentrated Extract Volume: 1000 (µL) Date Analyzed: 04/01/04 Time: 1730
 Soil Aliquot Volume: _____ (µL) Dilution Factor: 1 Analyst: TLS
 Injection Volume: 1 (µL) Prep Method: _____
 GPC Cleanup: (Y/N) N pH: _____ Analytical Method: OLMO 4.2
 Prep Batch: 271349 Analytical Batch: 271697 Sulfur Cleanup: (Y/N) N Instrument ID: GCS6A
 CONCENTRATION UNITS: ug/L Lab File ID: 2040331/SV6042

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|------------|---------------------|--------|---|---------|-------|
| 72-54-8 | 4,4'-DDD | 0.100 | U | 0.00010 | 0.100 |
| 72-55-9 | 4,4'-DDE | 0.100 | U | 0.00010 | 0.100 |
| 50-29-3 | 4,4'-DDT | 0.100 | U | 0.00010 | 0.100 |
| 309-00-2 | Aldrin | 0.050 | U | 0.00010 | 0.050 |
| 12674-11-2 | Aroclor-1016 | 1.00 | U | 0.00010 | 1.00 |
| 11104-28-2 | Aroclor-1221 | 2.00 | U | 0.00010 | 2.00 |
| 11141-16-5 | Aroclor-1232 | 1.00 | U | 0.00010 | 1.00 |
| 53469-21-9 | Aroclor-1242 | 1.00 | U | 0.00010 | 1.00 |
| 12672-29-6 | Aroclor-1248 | 1.00 | U | 0.00010 | 1.00 |
| 11097-69-1 | Aroclor-1254 | 1.00 | U | 0.00010 | 1.00 |
| 11096-82-5 | Aroclor-1260 | 1.00 | U | 0.00010 | 1.00 |
| 60-57-1 | Dieldrin | 0.100 | U | 0.00010 | 0.100 |
| 959-98-8 | Endosulfan I | 0.050 | U | 0.00010 | 0.050 |
| 33213-65-9 | Endosulfan II | 0.100 | U | 0.00010 | 0.100 |
| 1031-07-8 | Endosulfan sulfate | 0.100 | U | 0.00010 | 0.100 |
| 72-20-8 | Endrin | 0.100 | U | 0.00010 | 0.100 |
| 7421-93-4 | Endrin aldehyde | 0.100 | U | 0.00010 | 0.100 |
| 53494-70-5 | Endrin ketone | 0.100 | U | 0.00010 | 0.100 |
| 76-44-8 | Heptachlor | 0.050 | U | 0.00010 | 0.050 |
| 1024-57-3 | Heptachlor epoxide | 0.050 | U | 0.00010 | 0.050 |
| 72-43-5 | Methoxychlor | 0.500 | U | 0.00010 | 0.500 |
| 8001-35-2 | Toxaphene | 5.00 | U | 0.00010 | 5.00 |
| 319-84-6 | alpha-BHC | 0.050 | U | 0.00010 | 0.050 |
| 5103-71-9 | alpha-Chlordane | 0.050 | U | 0.00010 | 0.050 |
| 319-85-7 | beta-BHC | 0.050 | U | 0.00010 | 0.050 |
| 319-86-8 | delta-BHC | 0.050 | U | 0.00010 | 0.050 |
| 58-89-9 | gamma-BHC (Lindane) | 0.050 | U | 0.00010 | 0.050 |
| 5103-74-2 | gamma-Chlordane | 0.050 | U | 0.00010 | 0.050 |

ORGANICS ANALYSIS DATA SHEET

Lab Name: GCAL Sample ID: SKSW511009
 Lab Code: LA024 Case No.: _____ Contract: _____
 Matrix: Water SAS No: _____ SDG No.: 204032408
 Sample wt/vol: 1000 Units: mL Lab Sample ID: 20403240802
 Level: (low/mid) _____ Date Collected: 03/22/04 Time: 1040
 % Moisture: _____ decanted: (Y/N) _____ Date Received: 03/24/04
 GC Column: DB-608-30M ID: 53 (mm) Date Extracted: 3/26/04
 Concentrated Extract Volume: 1000 (µL) Date Analyzed: 04/02/04 Time: 1158
 Soil Aliquot Volume: _____ (µL) Dilution Factor: 1 Analyst: HJL
 Injection Volume: 1 (µL) Prep Method: _____
 GPC Cleanup: (Y/N) N pH: _____ Analytical Method: OLMO 4.2
 Prep Batch: 271349 Analytical Batch: 271697 Sulfur Cleanup: (Y/N) N Instrument ID: GCS6A

CONCENTRATION UNITS: µg/L

Lab File ID: 2040331/SV6053

CAS NO. COMPOUND

RESULT Q MDL RL

| | | | | | |
|------------|---------------------|-------|---|---------|-------|
| 72-54-8 | 4,4'-DDD | 0.100 | U | 0.00010 | 0.100 |
| 72-55-9 | 4,4'-DDE | 0.100 | U | 0.00010 | 0.100 |
| 50-29-3 | 4,4'-DDT | 0.100 | U | 0.00010 | 0.100 |
| 308-00-2 | Aldrin | 0.050 | U | 0.00010 | 0.050 |
| 12674-11-2 | Aroclor-1016 | 1.00 | U | 0.00010 | 1.00 |
| 11104-28-2 | Aroclor-1221 | 2.00 | U | 0.00010 | 2.00 |
| 11141-16-5 | Aroclor-1232 | 1.00 | U | 0.00010 | 1.00 |
| 53489-21-9 | Aroclor-1242 | 1.00 | U | 0.00010 | 1.00 |
| 12672-29-6 | Aroclor-1248 | 1.00 | U | 0.00010 | 1.00 |
| 11087-89-1 | Aroclor-1254 | 1.00 | U | 0.00010 | 1.00 |
| 11086-82-5 | Aroclor-1260 | 1.00 | U | 0.00010 | 1.00 |
| 80-57-1 | Dieldrin | 0.100 | U | 0.00010 | 0.100 |
| 959-98-8 | Endosulfan I | 0.050 | U | 0.00010 | 0.050 |
| 33213-65-9 | Endosulfan II | 0.100 | U | 0.00010 | 0.100 |
| 1031-07-8 | Endosulfan sulfate | 0.100 | U | 0.00010 | 0.100 |
| 72-20-8 | Endrin | 0.100 | U | 0.00010 | 0.100 |
| 7421-83-4 | Endrin aldehyde | 0.100 | U | 0.00010 | 0.100 |
| 53484-70-5 | Endrin ketone | 0.100 | U | 0.00010 | 0.100 |
| 76-44-8 | Heptachlor | 0.050 | U | 0.00010 | 0.050 |
| 1024-57-3 | Heptachlor epoxide | 0.050 | U | 0.00010 | 0.050 |
| 72-43-5 | Methoxychlor | 0.500 | U | 0.00010 | 0.500 |
| 8001-35-2 | Toxaphene | 5.00 | U | 0.00010 | 5.00 |
| 319-84-6 | alpha-BHC | 0.050 | U | 0.00010 | 0.050 |
| 5103-71-9 | alpha-Chlordane | 0.050 | U | 0.00010 | 0.050 |
| 319-85-7 | beta-BHC | 0.050 | U | 0.00010 | 0.050 |
| 319-86-8 | delta-BHC | 0.050 | U | 0.00010 | 0.050 |
| 58-89-9 | gamma-BHC (Lindane) | 0.050 | U | 0.00010 | 0.050 |
| 5103-74-2 | gamma-Chlordane | 0.050 | U | 0.00010 | 0.050 |

5/26/04
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ORGANICS ANALYSIS DATA SHEET

Lab Name: GCAL Sample ID: SKSW521009
 Lab Code: LA024 Case No.: _____ Contract: _____
 Matrix: Water SAS No.: _____ SDG No.: 204032408
 Sample wt/vol: 1000 Units: mL Lab Sample ID: 20403240806
 Level: (low/med) _____ Date Collected: 03/22/04 Time: 1230
 % Moisture: _____ decanted: (Y/N) _____ Date Received: 03/24/04
 GC Column: DB-608-30M ID: .53 (mm) Date Extracted: 3/26/04
 Concentrated Extract Volume: 1000 (µL) Date Analyzed: 04/01/04 Time: 1907
 Soil Aliquot Volume: _____ (µL) Dilution Factor: 1 Analyst: HJL
 Injection Volume: 1 (µL) Prep Method: _____
 GPC Cleanup: (Y/N) N pH: _____ Analytical Method: OLMO 4.2
 Prep Batch: 271349 Analytical Batch: 271697 Sulfur Cleanup: (Y/N) N Instrument ID: GCS6A
 Lab File ID: 2040331/SV6046

CONCENTRATION UNITS: ug/L

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|------------|---------------------|--------|---|---------|-------|
| 72-54-8 | 4,4'-DDD | 0.100 | U | 0.00010 | 0.100 |
| 72-55-9 | 4,4'-DDE | 0.100 | U | 0.00010 | 0.100 |
| 50-29-3 | 4,4'-DDT | 0.100 | U | 0.00010 | 0.100 |
| 309-00-2 | Aldrin | 0.050 | U | 0.00010 | 0.050 |
| 12674-11-2 | Aroclor-1016 | 1.00 | U | 0.00010 | 1.00 |
| 11104-28-2 | Aroclor-1221 | 2.00 | U | 0.00010 | 2.00 |
| 11141-16-5 | Aroclor-1232 | 1.00 | U | 0.00010 | 1.00 |
| 53469-21-9 | Aroclor-1242 | 1.00 | U | 0.00010 | 1.00 |
| 12672-29-6 | Aroclor-1248 | 1.00 | U | 0.00010 | 1.00 |
| 11097-69-1 | Aroclor-1254 | 1.00 | U | 0.00010 | 1.00 |
| 11096-82-5 | Aroclor-1260 | 1.00 | U | 0.00010 | 1.00 |
| 60-57-1 | Dieldrin | 0.100 | U | 0.00010 | 0.100 |
| 959-98-8 | Endosulfan I | 0.050 | U | 0.00010 | 0.050 |
| 33213-65-9 | Endosulfan II | 0.100 | U | 0.00010 | 0.100 |
| 1031-07-8 | Endosulfan sulfate | 0.100 | U | 0.00010 | 0.100 |
| 72-20-8 | Endrin | 0.100 | U | 0.00010 | 0.100 |
| 7421-93-4 | Endrin aldehyde | 0.100 | U | 0.00010 | 0.100 |
| 53494-70-5 | Endrin ketone | 0.100 | U | 0.00010 | 0.100 |
| 76-44-8 | Heptachlor | 0.050 | U | 0.00010 | 0.050 |
| 1024-57-3 | Heptachlor epoxide | 0.050 | U | 0.00010 | 0.050 |
| 72-43-5 | Methoxychlor | 0.500 | U | 0.00010 | 0.500 |
| 8001-35-2 | Toxaphene | 5.00 | U | 0.00010 | 5.00 |
| 319-84-6 | alpha-BHC | 0.050 | U | 0.00010 | 0.050 |
| 5103-71-9 | alpha-Chlordane | 0.050 | U | 0.00010 | 0.050 |
| 319-85-7 | beta-BHC | 0.050 | U | 0.00010 | 0.050 |
| 319-86-8 | delta-BHC | 0.050 | U | 0.00010 | 0.050 |
| 58-89-9 | gamma-BHC (Lindane) | 0.050 | U | 0.00010 | 0.050 |
| 5103-74-2 | gamma-Chlordane | 0.050 | U | 0.00010 | 0.050 |

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INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKSWE81009

Lab Name: GCAL Contract: _____
Lab Code: LA024 Case No. _____ SAS No. _____ SDG No.: 204032408
Matrix (soil / water): Water Lab Sample ID: 20403240801
Level: (low / med) _____ Date Received: 03/24/04
% Solids: _____

Concentration Units (ug/L or mg/kg dry weight): ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|-----|----|
| 7440-38-0 | Antimony | 3.7 | U | N | P |
| 7440-38-2 | Arsenic | 2.9 | U | N,* | P |
| 7440-39-3 | Barium | 0.3 | B | *E | P |
| 7440-41-7 | Beryllium | 0.1 | U | | P |
| 7440-43-0 | Cadmium | 0.2 | U | | P |
| 7440-47-3 | Chromium | 1.3 | B | *E | P |
| 7440-50-8 | Copper | 6.5 | B | * | P |
| 7439-89-6 | Iron | 37.8 | B | *E | P |
| 7439-92-1 | Lead | 1.5 | U | * | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 1.4 | B | *E | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 1.2 | B | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-66-6 | Zinc | 9.7 | B | N,* | P |
| 57-12-5 | Cyanide | 0.5 | U | | AS |

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Color Before: COLORLESS Clarity Before: CLEAR Texture: _____
Color After: COLORLESS Clarity After: CLEAR Artifacts: _____
Comments: _____

EPA SAMPLE NO.

SKSW511009

Lab Name: GCAL Contract:
 Lab Code: LA024 Case No.: SAS No.: SDG No.: 204032408
 Matrix: (soil / water) Water Lab Sample ID: 20403240802
 Level: (low / med) Date Received: 03/24/04
 % Solids:

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|-----|----|
| 7440-36-0 | Antimony | 7.3 | B | N | P |
| 7440-38-2 | Arsenic | 50.7 | | N,* | P |
| 7440-39-3 | Barium | 499 | | *,E | P |
| 7440-41-7 | Beryllium | 4.9 | B | | P |
| 7440-43-9 | Cadmium | 5.0 | | | P |
| 7440-47-3 | Chromium | 72.6 | | *,E | P |
| 7440-50-8 | Copper | 131 | | * | P |
| 7439-89-6 | Iron | 124000 | | *,E | P |
| 7439-92-1 | Lead | 122 | | * | P |
| 7439-97-6 | Mercury | 0.1 | B | | AV |
| 7440-02-0 | Nickel | 116 | | *,E | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-66-6 | Zinc | 490 | | N,* | P |
| 57-12-5 | Cyanide | 0.8 | B | | AS |

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Color Before: DK. BROWN

Clarity Before: CLOUDY

Texture:

Color After: LT. BROWN

Clarity After: **CLEAR**

Artifacts:

Comments:

EPA SAMPLE NO.

SKSW51DUP1009

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|-------------------------|-------|---------------|--------------------|
| Lab Name: | GCAL | Contract | |
| Lab Code: | LA024 | Case No. | |
| | | SAS No. | SDG No.: 204032408 |
| Matrix (soil / water) | Water | Lab Sample ID | 20403240805 |
| Level: (low / med) | | Date Received | 03/24/04 |
| % Solids: | | | |

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|-----|----|
| 7440-38-0 | Antimony | 6.7 | B | N | P |
| 7440-38-2 | Arsenic | 67.0 | | N,* | P |
| 7440-39-3 | Barium | 714 | | *E | P |
| 7440-41-7 | Beryllium | 6.7 | | | P |
| 7440-43-9 | Cadmium | 7.2 | | | P |
| 7440-47-3 | Chromium | 102 | | *E | P |
| 7440-50-8 | Copper | 174 | | * | P |
| 7439-89-6 | Iron | 173000 | | *E | P |
| 7439-92-1 | Lead | 169 | | * | P |
| 7439-97-6 | Mercury | 0.3 | | | AV |
| 7440-02-0 | Nickel | 162 | | *E | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-66-6 | Zinc | 679 | | N,* | P |
| 57-12-5 | Cyanide | 1.1 | B | | AS |

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|----------------------|------------------|------------------------|---------------|-------------------|-------------------|
| Color Before: | <u>DK. BROWN</u> | Clarity Before: | <u>CLOUDY</u> | Texture: | <u> </u> |
| Color After: | <u>LT. BROWN</u> | Clarity After: | <u>CLEAR</u> | Artifacts: | <u> </u> |
| Comments: | | | | | |

EPA SAMPLE NO.

SKSW521009

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|-----|----|
| 7440-36-0 | Antimony | 3.7 | U | N | P |
| 7440-38-2 | Arsenic | 11.1 | | N,* | P |
| 7440-39-3 | Barium | 112 | B | *,E | P |
| 7440-41-7 | Beryllium | 0.7 | B | | P |
| 7440-43-9 | Cadmium | 0.5 | B | | P |
| 7440-47-3 | Chromium | 12.7 | | *,E | P |
| 7440-50-8 | Copper | 22.0 | B | * | P |
| 7439-89-6 | Iron | 17800 | | *,E | P |
| 7439-92-1 | Lead | 17.3 | | * | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 16.4 | B | *,E | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-66-6 | Zinc | 52.9 | | N,* | P |
| 57-12-5 | Cyanide | 0.6 | B | | AS |

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U.S. EPA - CLP
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKSW531009

Lab Name: GCAL Contract: _____
Lab Code: LA024 Case No: _____ SAS No: _____ SDG No.: 204032408
Matrix: (soil / water) Water Lab Sample ID: 20403240807
Level: (low / med) _____ Date Received: 03/24/04
% Solids: _____

Concentration Units (ug/L or mg/kg dry weight) ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|-----|----|
| 7440-38-0 | Antimony | 37 | U | N | P |
| 7440-38-2 | Arsenic | 29 | U | N,* | P |
| 7440-39-3 | Barium | 417 | B | *E | P |
| 7440-41-7 | Beryllium | 01 | U | | P |
| 7440-43-9 | Cadmium | 02 | U | | P |
| 7440-47-3 | Chromium | 17 | B | *E | P |
| 7440-50-8 | Copper | 32 | B | * | P |
| 7439-89-6 | Iron | 67.8 | B | *E | P |
| 7439-92-1 | Lead | 1.5 | U | * | P |
| 7439-97-8 | Mercury | 01 | U | | AV |
| 7440-02-0 | Nickel | 1.1 | B | *E | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 06 | B | | P |
| 7440-28-0 | Thallium | 26 | U | | P |
| 7440-66-6 | Zinc | 06 | U | N,* | P |
| 57-12-5 | Cyanide | 05 | U | | AS |

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Color Before: COLORLESS Clarity Before: CLEAR Texture: _____
Color After: COLORLESS Clarity After: CLEAR Artifacts: _____
Comments: _____

U.S. EPA - CLP
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INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKSWE1009 (DISS)

Lab Name: GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204032408
Matrix: (soil / water) Water Lab Sample ID: 20403240808
Level: (low / med) _____ Date Received: 03/24/04
% Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7440-36-0 | Antimony | 3.7 | U | | P |
| 7440-38-2 | Arsenic | 2.9 | U | N | P |
| 7440-39-3 | Barium | 0.4 | B | | P |
| 7440-41-7 | Beryllium | 0.1 | U | | P |
| 7440-43-9 | Cadmium | 0.2 | U | | P |
| 7440-47-3 | Chromium | 0.8 | U | | P |
| 7440-50-8 | Copper | 4.2 | B | | P |
| 7439-89-6 | Iron | 14.1 | U | | P |
| 7439-92-1 | Lead | 1.5 | U | | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 0.7 | U | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-66-6 | Zinc | 5.6 | B | | P |

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Color Before: COLORLESS Clarity Before: CLEAR Texture: _____
Color After: COLORLESS Clarity After: CLEAR Artifacts: _____
Comments: _____

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INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKSW511009 (DISS)

Lab Name: GCAL Contract: _____
 Lab Code: LA024 Case No: _____ SAS No: _____ SDG No.: 204032408
 Matrix (soil / water): Water Lab Sample ID: 20403240809
 Level (low / med): _____ Date Received: 03/24/04
 % Solids: _____

Concentration Units (ug/L or mg/kg dry weight): ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7440-36-0 | Antimony | 37 | U | | P |
| 7440-38-2 | Arsenic | 29 | U | N | P |
| 7440-39-3 | Barium | 35.8 | B | | P |
| 7440-41-7 | Beryllium | 01 | U | | P |
| 7440-43-9 | Cadmium | 02 | U | | P |
| 7440-47-3 | Chromium | 16 | B | | P |
| 7440-50-8 | Copper | 43 | B | | P |
| 7439-89-6 | Iron | 14.1 | U | | P |
| 7439-92-1 | Lead | 15 | U | | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 1.0 | B | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-66-6 | Zinc | 0.6 | U | | P |

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Color Before: COLORLESS Clarity Before: CLEAR Texture: _____
 Color After: COLORLESS Clarity After: CLEAR Artifacts: _____
 Comments: _____

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INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKSW51DUP1009 (DISS)

Lab Name: GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204032408
Matrix: (soil / water) Water Lab Sample ID: 20403240811
Level: (low / med) _____ Date Received: 03/24/04
% Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7440-36-0 | Antimony | 3.7 | U | | P |
| 7440-38-2 | Arsenic | 2.9 | U | N | P |
| 7440-39-3 | Barium | 33.0 | B | | P |
| 7440-41-7 | Beryllium | 0.1 | U | | P |
| 7440-43-9 | Cadmium | 0.2 | U | | P |
| 7440-47-3 | Chromium | 1.4 | B | | P |
| 7440-50-8 | Copper | 1.6 | B | | P |
| 7439-89-6 | Iron | 14.1 | U | | P |
| 7439-92-1 | Lead | 1.5 | U | | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 1.0 | B | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-66-6 | Zinc | 0.6 | U | | P |

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Color Before: COLORLESS Clarity Before: CLEAR Texture: _____
Color After: COLORLESS Clarity After: CLEAR Artifacts: _____
Comments: _____

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INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKSW521009 (DISS)

Lab Name: GCAL Contract: _____
 Lab Code: LA024 Case No: _____ SAS No: _____ SDG No.: 204032408
 Matrix: (soil / water) Water Lab Sample ID: 20403240812
 Level: (low / med) _____ Date Received: 03/24/04
 % Solids: _____
 Concentration Units (ug/L or mg/kg dry weight): ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7440-38-0 | Antimony | 3.7 | L | | P |
| 7440-38-2 | Arsenic | 2.9 | L | N | P |
| 7440-39-3 | Barium | 36.7 | B | | P |
| 7440-41-7 | Beryllium | 0.1 | L | | P |
| 7440-43-9 | Cadmium | 0.2 | L | | P |
| 7440-47-3 | Chromium | 1.7 | B | | P |
| 7440-50-8 | Copper | 3.0 | B | | P |
| 7439-89-6 | Iron | 14.1 | U | | P |
| 7439-92-1 | Lead | 1.5 | U | | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 2.4 | B | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | L | | P |
| 7440-28-0 | Thallium | 2.6 | L | | P |
| 7440-66-6 | Zinc | 3.4 | B | | P |

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Color Before: COLORLESS Clarity Before: CLEAR Texture: _____
 Color After: COLORLESS Clarity After: CLEAR Artifacts: _____
 Comments: _____

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INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKSW531009 (DISS)

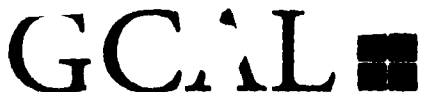
Lab Name: GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204032408
Matrix: (soil / water) Water Lab Sample ID: 20403240813
Level: (low / med) _____ Date Received: 03/24/04
% Solids: _____
Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7440-36-0 | Antimony | 3.7 | U | | P |
| 7440-38-2 | Arsenic | 2.9 | U | N | P |
| 7440-39-3 | Barium | 40.7 | B | | P |
| 7440-41-7 | Beryllium | 0.1 | U | | P |
| 7440-43-9 | Cadmium | 0.2 | U | | P |
| 7440-47-3 | Chromium | 1.3 | B | | P |
| 7440-50-8 | Copper | 2.0 | B | | P |
| 7439-89-6 | Iron | 14.1 | U | | P |
| 7439-92-1 | Lead | 1.5 | U | | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 0.7 | U | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-66-6 | Zinc | 4.5 | B | | P |

R

6/3/09
msc

Color Before: COLORLESS Clarity Before: CLEAR Texture: _____
Color After: COLORLESS Clarity After: CLEAR Artifacts: _____
Comments: _____



GULF COAST ANALYTICAL LABORATORIES, INC.
7979 GSRI Avenue, Baton Rouge, Louisiana 70820-7402
Phone 225.789.4900 • Fax 225.767.5717

CHAIN OF CUSTODY RECORD

Lab use only

Earth Tech

Client Name

4342

Client #

284032468

Workorder #

4/17/04

Due Date

Report to:

Client: Earth Tech
Address: 200 Vine Street
Wilder, LA 70776
Contact: Pat Higgins
Phone: 859 442-2300
Fax: 859 442-2311

Bill to:

Client: SAMS
Address: SAMS
Contact: SAMS
Phone: SAMS
Fax: SAMS

Analytical Requests & Method

Lab use only:

Custody Seal

used ☒ yes ☐ no

in tact ☒ yes ☐ no

Temperature °C

P.O. Number

54280

Project Name/Number

Chinner Landfill - 1 Qtr. 2004

Sampled By:

Pat Higgins/Chris Cox

| Matrix | Date | Time (400) | Sample Description | Preservatives | No. Containers | Volatiles | Semi-volatiles | Pesticides | PCBs | Trace Metals | Dissolved Metals | Cyanide | Other |
|--------|---------|------------|---------------------|---------------|----------------|-----------|----------------|------------|------|--------------|------------------|---------|-------|
| W | 3/24/04 | 1115 | X SK SW SI MSD 1009 | Various | 10 | X | X | X | X | X | X | X | MSD 4 |
| W | 3/24/04 | 1315 | X SK SW ED 1009 | Various | 10 | X | X | X | X | X | X | X | 8 |
| W | 3/24/04 | 1040 | X SK SW SI 1009 | HCL | 3 | X | | | | | | | |
| W | 3/24/04 | 1055 | X SK SW SI MS 1009 | HCL | 3 | X | | | | | | | |
| W | 3/24/04 | 1210 | X SK SW SI 1009 | HCL | 3 | X | | | | | | | |
| W | 3/24/04 | 1255 | X SK SW SI 1009 | HCL | 3 | X | | | | | | | |

Remarks:

Refer to Chap. 5
Table 7
(TEL) and
Table 8
(TAL) of
The Final
O&M Plan
for the
list of
analytes

Turn Around Time: ☐ 24-48 hrs. ☐ 3 days ☐ 1 week ☒ Standard ☐ Other

Relinquished by: (Signature)

Pat Higgins

Relinquished by: (Signature)

FED EX # 12549688324

Relinquished by: (Signature)

Pat Higgins

Received by: (Signature)

FED EX

Received by: (Signature)

Chris Cox

Received by: (Signature)

Chris Cox

Date:

3/23/04

Date:

3/24/04

Date:

3/24/04

Time:

0920

Time:

0920

Time:

0920

Note:

- No trip blank submitted due to bottles breaking during shipment
- Low sample volume for Equipment Blank

By submitting these samples, you agree to the terms and conditions contained in our most recent schedule.



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Phone 225.769.4900 • Fax 225.767.5717

CHAIN OF CUSTODY RECORD

Lab use only

Earth Tech

Client Name

4342

Client #

204033408

Workorder #

4/7/04

Due Date

Report to:

Client: Earth Tech
Address: 200 Vine Street
Wilders, KY 40376
Contact: Pat Higgins
Phone: 859 442-2300
Fax: 859 442-2311

Bill to:

Client: _____
Address: _____
Contact: SAMA
Phone: _____
Fax: _____

Analytical Requests & Method

| Volatiles | Semi-volatiles | Pesticides | PCB's | Total Metals | Dissolved Metals | Cyanide | Diss |
|-----------|----------------|------------|-------|--------------|------------------|---------|------|
| | X | X | X | X | X | X | 9 |
| | X | X | X | X | X | X | 10 |

Lab use only:

Custody Seal
used ☒ yes ☐ no
in tact ☒ yes ☐ no
Temperature °C 3

P.O. Number

54280

Project Name/Number

Shinner Landfill- 1 Qtr. 2004

Sampled By:

Pat Higgins / Chris Cox

| Matrix | Date | Time (2400) | Comp | Grab | Sample Description | Preservatives | No Containers |
|--------|---------|-------------|------|------|--------------------|---------------|---------------|
| W | 3/23/04 | 1040 | | X | Sk SW 51 1009 | Various | 7 |
| W | 3/23/04 | 1055 | | X | Sk SW 51 MS 1009 | Various | |

Lab ID

3/24

Remarks:

Refer to
Table 7
(TCL) and
Table 8
(TAC) of
the final
O&M Plan
for the
list of
analytes

Turn Around Time: ☐ 24-48 hrs. ☐ 3 days ☐ 1 week ☒ Standard ☐ Other

Relinquished by: (Signature)

Received by: (Signature)

Date: 3/23/04 Time:

Note:

Relinquished by: (Signature)

Received by: (Signature)

Date: 3/24/04 Time: 1520

Relinquished by: (Signature)

Received by: (Signature)

Date: Time:

By submitting these samples, you agree to the terms and conditions contained in our most recent schedule of services.



GULF COAST ANALYTICAL LABORATORIES, INC.
7979 GSRI Avenue, Baton Rouge, Louisiana 70820-7402
Phone 225 789 4900 • Fax 225 787 5717

CHAIN OF CUSTODY

RECORD

Lab use only

Earth Tech
Client Name

4342
Client #

704032408
Workorder #

4/7/04
Due Date

Report to:

Bill to:

Analytical Requests & Method

Lab use only.

Client: Earth Tech
Address: 200 Vine Street
Wilder, KY 41076
Contact: Pat Higgins
Phone: 859 442-2300
Fax: 859 442-2311

Client: Earth Tech
Address: Earth Tech
Contact: Earth Tech
Phone: Earth Tech
Fax: Earth Tech

Custody Seal

used ☒ yes ☐ no

in tact ☒ yes ☐ no

Temperature °C

30

P.O. Number

54280

Project Name/Number

Shinner Landfill - 1 Qtr. 2004

Sampled By:

Pat Higgins / Chris Cox

| Matrix | Date | Time (2400) | Sample Description | Preservatives | No Containers |
|--------|---------|-------------|--------------------|---------------|---------------|
| W | 3/24/04 | 12:10 | X Sk SW 52 1009 | Various | 7 |
| W | 3/24/04 | 12:55 | X Sh SW 53 1009 | Various | 7 |

Volatiles
Semi-Volatiles
Pesticides
PCBs
total metals
Dissolved metals
Cyanide
455
12
13

Remarks:

Refer to
Table 7
(TCL) and
Table 8
(TAL) of
the Final
O&M Plan
for the
list of
analytes

3/24
-6
1

Turn Around Time: ☐ 24-48 hrs. ☐ 3 days ☐ 1 week ☒ Standard ☐ Other

Relinquished by: (Signature)
Pat Higgins
Relinquished by: (Signature)
Chris Cox
Relinquished by: (Signature)
Chris Cox

Received by: (Signature)
FED EX
Received by: (Signature)
Chris Cox
Received by: (Signature)
Chris Cox

Date: 3/23/04 Time:
Date: 3/24/04 Time: 0920
Date: Time:

Note:

By submitting these samples, you agree to the terms and conditions contained in our most recent schedule of services.

DATA VALIDATION REPORT
FOR
SKINNER LANDFILL SITE
EARTH TECH: PROJECT NUMBER 54280
LABORATORY REPORT NUMBER 204031909
PROJECT MANAGER: Ron Rolker
Date: May 11, 2004
Data Validator: Mark Kromis

APPENDIX C
LIST OF ACRONYMS

| | |
|---------------|---|
| BFB | Bromofluorobenzene |
| CC | Continuing Calibration |
| CCV | Continuing Calibration Verification |
| CCB | Continuing Calibration Blanks |
| CLP | Contract Laboratory Program |
| CRDL | Contract Required Detection Limit |
| DFTPP | Decafluorotriphenylphosphine |
| GC/MS | Gas Chromatograph Mass Spectrometer |
| IC | Initial Calibration |
| ICB | Initial Calibration Blank |
| IDL | Instrument Detection Limit |
| ICP | Inductively Coupled Plasma |
| ICS | Interference Check Sample |
| ICV | Initial Calibration Verification |
| ILM | Inorganic Analysis Multi-Media Multi-Concentration |
| INDAM | Individual A Mixture |
| INDBM | Individual B Mixture |
| mg/L | milligrams per liter |
| MS/MSD | Matrix Spike Matrix Spike Duplicate |
| OLC | Organic Analysis Low Concentration |
| OLM | Organic Analysis Multi-Media Multi-Concentration |
| %D | Percent Difference |
| % RSD | Percent Relative Standard Deviation |
| PB | Preparation Blanks |
| QC | Quality Control |
| RF | Response Factor |
| RPD | Relative Percent Difference |
| RRF | Relative Response Factor |
| SDG | Sample Delivery Group |
| SOW | Statement of Work |
| µg/L | micrograms per liter |
| US EPA | United States Environmental Protection Agency |
| VOC | Volatile Organic Compounds |
| VTSR | Validated Time of Sample Receipt |

DATA VALIDATION SUMMARY – SAMPLE DELIVERY GROUP 204031909 INORGANICS

Validation of the inorganics data, as prepared by Gulf Coast Analytical Laboratories (GCAL) for the samples collected from the Skinner Landfill site in March 2004, was conducted by Earth Tech using the National Functional Guidelines for Inorganic Data Review, (US EPA, February, 1994), as appropriate. The results were reported by GCAL under Sample Delivery Group (SDG) 204031909.

| GCAL # | Sample Description |
|-------------|----------------------|
| 20403190901 | SKGW581009 |
| 20403190902 | SKGW58DUP1009 |
| 20403190903 | SKGW631009 |
| 20403190904 | SKGW611009 |
| 20403190905 | SKGW61MS1009 |
| 20403190907 | SKGW61DUP1009 |
| 20403190908 | SKGWFB1009 |
| 20403190910 | SKGW581009 (DISS) |
| 20403190911 | SKGW58DUP1009 (DISS) |
| 20403190912 | SKGW631009 (DISS) |
| 20403190913 | SKGW611009 (DISS) |
| 20403190914 | SKGW61MS1009 (DISS) |
| 20403190915 | SKGW61DUP1009 (DISS) |
| 20403190916 | SKGWFB1009 (DISS) |
| 20403220801 | SKGW06R1009 |
| 20403220802 | SKGW07R1009 |
| 20403220803 | SKGW591009 |
| 20403220804 | SKGW601009 |
| 20403220805 | SKGW62A1009 |
| 20403220806 | SKGW641009 |
| 20403220807 | SKGW651009 |
| 20403220808 | SKGW06R1009 (DISS) |
| 20403220809 | SKGW07R1009 (DISS) |
| 20403220810 | SKGW591009 (DISS) |
| 20403220811 | SKGW601009 (DISS) |
| 20403220812 | SKGW62A1009 (DISS) |
| 20403220813 | SKGW641009 (DISS) |

INTRODUCTION

Analyses of metals were performed according to Contract Laboratory Program (CLP)-Inorganic Analysis Multi-media Multi-concentration ILM04.1 Statement of Work (SOW). Results of the sample analyses are reported by the laboratory as either qualified or unqualified. Unqualified results mean that the reported values maybe used without reservation. Various qualifier codes are used by the laboratory to denote specific information regarding the analytical results.

The data validation process is intended to evaluate the data on a technical basis. The data package also was subjected to an internal laboratory quality review prior to submission to Earth Tech for data validation.

During the validation process, laboratory-qualified and unqualified data are verified against all available supporting documentation. Based on this evaluation, qualifier codes may be added, deleted or modified by the data user. Final results are therefore, either qualified or unqualified. Validator-qualified results are annotated with the following codes in accordance with the Functional Guidelines:

- U** The constituent was analyzed for, but was not detected above the level of the associated analytical reporting limit. The associated value is either the sample quantitation limit or the sample detection limit.
- J** The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample.
- UJ** The analyte was not detected above the reported sample quantitation limit. However, the reported quantitation limit is approximate and may or may not represent the actual limit of quantitation necessary to accurately and precisely measure the analyte in the sample.
- R** The sample results are rejected due to serious deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte cannot be verified.

Details of the inorganics data validation findings and conclusions are provided in the following sections of this report:

1. Holding Times
2. Calibration
 - A. Initial Calibration (IC)
 - B. Continuing Calibration (CC)
3. Blanks
4. Inductively Coupled Plasma (ICP) Interference Check Sample
5. Laboratory Control Sample (LCS)
6. Duplicate Analysis

7. Spike Sample Analysis
8. ICP Serial Dilution
9. System Performance
10. Documentation
11. Overall Assessment

1. HOLDING TIMES

All samples for inorganics analyses were analyzed within the 180-day holding time for preserved aqueous samples. Mercury analyses were conducted within the 28-day holding time for aqueous samples undergoing CLP protocol. Cyanide analyses were conducted within the 14-day holding time. The cooler temperature upon receipt at the laboratory was within the recommended temperature of 4°C +/- 2°C.

2. CALIBRATION

A. Initial Calibration

The percent recoveries for the Initial Calibration Verification (ICV) standard were within Quality Control (QC) limits for all constituents.

B. Continuing Calibration

The percent recoveries for the Continuing Calibration Verification (CCV) standard were within QC limits for all constituents.

3. BLANKS

The Initial Calibration Blank (ICB), Continuing Calibration Blanks (CCB) and Preparation Blanks (PB) were analyzed at the appropriate frequencies. No analytes were detected in the ICB, CCB, PB and Field blanks above the corresponding Contract Required Detection Limit (CRDL).

4. ICP INTERFERENCE CHECK SAMPLE

Results for the ICP analysis of the Interference Check Sample (ICS) solution AB were within 20% of the true value.

5. LABORATORY CONTROL SAMPLES

Recoveries were within the control limit (80-120%) for all constituents.

6. DUPLICATE ANALYSIS

The Relative Percent Difference (RPD) between the sample and duplicate results were within the acceptance criteria for all target compounds.

7. SPIKE SAMPLE ANALYSIS

The laboratory used sample SKGW611009 for the matrix spike sample. The MS percent recoveries were within the acceptance criteria (75%-125%) with the exception of Selenium in the total (45%) and dissolved (73%) fractions. As per the National Functional Guidelines: if the percent recovery is greater than 30% and less than 74% qualify detected results for that analyte with "J" and non-detected results with "UJ".

8. ICP SERIAL DILUTION

As noted in the National Functional Guidelines: If the analyte concentration is at least 50 times above the IDL, its serial dilution analysis must then agree within 10% of the original determination after corrected for dilution. The serial dilution is performed to determine whether any significant chemical or physical interference's exist due to matrix effects. The serial dilution percent differences were within the acceptance criteria for all target analytes.

9. SYSTEM PERFORMANCE

The analytical system appears to have been working well at the time of these analyses, based on the evaluation of the raw data.

10. DOCUMENTATION

The documentation appeared accurate and in order.

11. OVERALL ASSESSMENT

The percent recoveries for Thallium in the Contract Required Detection Limit (CRDL) standards were 125.8%, 146.7%, and 130.4%. The percent recoveries for Zinc in the Contract Required Detection Limit (CRDL) standards were 73.9%, 71.5%, and 68.9%. The detected Thallium and Zinc results greater than the IDL but less than two times the CRDL were qualified with as estimated with "J". The percent recovery for Selenium in the last Contract Required Detection Limit (CRDL) standards was 138.5%. The Selenium results were previously qualified under Section 7-titled "Spike Sample Analysis". The results are acceptable with the validator-added qualifiers.

**DATA VALIDATION SUMMARY – SAMPLE DELIVERY GROUP 204031909
SEMIVOLATILE ORGANICS**

Validation of the Gas Chromatograph/Mass Spectrometer (GC/MS) semi-volatile organics data, as prepared by Gulf Coast Analytical Laboratories (GCAL) for the samples collected from the Skinner Landfill site in March 2004, was conducted by Earth Tech using the National Functional Guidelines for Organic Data Review, (US EPA, October, 1999) as appropriate. The results were reported by GCAL under SDG 204031909.

| GCAL # | Sample Description |
|---------------|---------------------------|
| 20403190901 | SKGW581009 |
| 20403190902 | SKGW58DUP1009 |
| 20403190903 | SKGW631009 |
| 20403190904 | SKGW611009 |
| 20403190905 | SKGW61MS1009 |
| 20403190906 | SKGW61MSD1009 |
| 20403190908 | SKGWFB1009 |
| 20403220801 | SKGW06R1009 |
| 20403220802 | SKGW07R1009 |
| 20403220803 | SKGW591009 |
| 20403220804 | SKGW601009 |
| 20403220805 | SKGW62A1009 |
| 20403220806 | SKGW641009 |
| 20403220807 | SKGW651009 |

INTRODUCTION

Analyses were performed according to CLP-Organic Analysis Multi-Media, Multi-Concentration OLM04.2 SOW. Results of the sample analyses are reported by the laboratory as either qualified or unqualified. Unqualified results mean that the reported values may be used without reservation. Various data qualifier codes are used by the laboratory to denote specific information regarding the analytical results. The data validation process is intended to evaluate the data on a technical basis. The data package also was subjected to an internal laboratory quality review prior to submission to Earth Tech for data validation.

During the validation process, laboratory-qualified and unqualified data are verified against all available supporting documentation. Based on this evaluation, qualifier codes may be added, deleted or modified by the data user. Final results are therefore, either qualified or unqualified. Validator-qualified results are annotated with the following codes in accordance with the Functional Guidelines:

- U The constituent was analyzed for, but was not detected above the level of the associated analytical reporting limit. The associated value is either the sample quantitation limit or the sample detection limit.

- J** The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample.
- UJ** The analyte was not detected above the reported sample quantitation limit. However, the reported quantitation limit is approximate and may or may not represent the actual limit of quantitation necessary to accurately and precisely measure the analyte in the sample.
- R** The sample results are rejected due to serious deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte cannot be verified.

Details of the semivolatile data validation findings and conclusions are provided in the following sections of this report:

- 1. Holding Times**
- 2. GC/MS Tuning**
- 3. Calibration**
 - A. IC**
 - B. CC**
- 4. Blanks**
- 5. System Monitoring Compound Recovery**
- 6. MS/MSD**
- 7. Internal Standards Performance**
- 8. Compound Identification**
- 9. Constituent Quantitation and Reported Detection Limits**
- 10. System Performance**
- 11. Documentation**
- 12. Overall Assessment**

1. HOLDING TIMES

All samples were initially extracted within the seven-day technical holding time and the five-day VTSR method holding time. The cooler temperature upon receipt at the laboratory was within the recommended temperature of 4°C +/- 2°C.

2. GC/MS TUNING

The samples were analyzed on a single GC/MS system, identified as MSSV2. Two decafluorotriphenylphosphine (DFTPP) tunes were run representing the shift in which the standards and samples were analyzed. The DFTPP tunes are acceptable.

3. CALIBRATION

A. Initial Calibration

One IC dated 3/19/04 was analyzed in support of the semivolatile sample analyses. Documentation of the IC was present in the data package, and the Relative Response Factor (RRF), as well as percent % RSD values were accurately reported for all target compounds. The criteria employed for technical data review purposes are different than those used in the method. The laboratory must meet a minimum RRF of 0.01; however, for data review purposes, a RRF criterion of "greater than or equal to 0.05" is applied to all semi-volatile compounds. The RRF's and the average RRF were within the acceptance criteria specified in the method for all reported analytes. The %RSD's were within the acceptance criteria specified in the method for all target analytes with the exception of Di-n-butylphthalate (32.0%), Di-n-octylphthalate (30.3%) and Diethylphthalate (41.8%). The lowest point of the calibration curve was dropped for Di-n-butylphthalate, Di-n-octylphthalate, and Diethylphthalate and the %RSD were recalculated. The recalculated %RSD were within the acceptance criteria of less than 30%. Di-n-octylphthalate and Diethylphthalate were not detected in the associated samples therefore data qualification was not required. The detected results for Di-n-butylphthalate were mitigated do to the presence of Di-n-butylphthalate in the associated method blank.

B. Continuing Calibration

One CC dated 3/26/04 was analyzed in support of the semivolatile sample analyses reported in the data submissions. The RRF's for the CC dated 3/26/04 were within the acceptance criteria. The percent difference (%D) between the average RRF's and the CC Response Factors for the CC dated 3/26/04 were within the acceptance criteria with the exception the %D for Di-n-butylphthalate and Di-n-octylphthalate. As per the National Functional Guidelines, if the %D exceeds the acceptance criteria qualify detected results for that analyte with "J" and non-detected results for that analyte with "UJ".

4. BLANKS

One laboratory semivolatile method blank and field blank were analyzed with this SDG. The results are summarized below.

Method Blank (0322SBLK)

Di-n-butylphthalate was detected at a concentration of 0.604 ppb in method blank 0322SBLK.

Field Blank (SKGWFB1009)

There were no analytes detected above the MDL in the field blank collected on 3/18/04.

5. SYSTEM MONITORING COMPOUND RECOVERY

All reported semivolatile system monitoring compounds were recovered within acceptable control limits.

6. MATRIX SPIKE/MATRIX SPIKE DUPLICATE (MS/MSD)

Sample SKGWD611009 was submitted for MS/MSD analysis. The MS/MSD percent recoveries were within the acceptance criteria with the exception of the 4-Nitrophenol. The %RPD between the MS/MSD are within the acceptance criteria. As per the National Functional Guidelines, no action is taken on MS/MSD data alone.

7. INTERNAL STANDARDS PERFORMANCE

Internal standard areas and retention times were within acceptable limits for the reported semivolatile sample analyses.

8. COMPOUND IDENTIFICATION

All reported semivolatile constituents were correctly identified with supporting chromatograms present in the data package.

9. CONSTITUENT QUANTITATION AND REPORTED DETECTION LIMITS

Constituent quantitations were correctly calculated and reported for semivolatile constituents.

10. SYSTEM PERFORMANCE

The analytical system appears to have been working well at the time of these analyses, based on the evaluation of the raw data submitted for review.

11. DOCUMENTATION

The documentation appeared accurate and in order with the exception of the following: GCAL inadvertently left the Date Extracted off of the CLP Form 1's therefore the data validator inserted the Date Extracted on the CLP Form 1's. GCAL also inadvertently left the "B" qualifier off of the CLP Form 1's for the compound Di-n-butylphthalate therefore the data validator inserted a "B" qualifier in the "Q" column of the CLP Form 1's.

The "B" qualifier indicates that the analyte was detected in the associated method blank. On pages 210/212/213/214 sample SKGW07R1009 was not listed but 2 field blanks were listed while only one field blank submitted for analysis. GCAL corrected the mistake and reissued corrected pages.

12. OVERALL ASSESSMENT

The results are acceptable with the validator-added qualifiers.

DATA VALIDATION SUMMARY – SAMPLE DELIVERY GROUP 204031909 VOLATILE ORGANIC

Validation of the GC/MS volatile organics data, as prepared by Gulf Coast Analytical Laboratories (GCAL) for the samples collected from the Skinner Landfill site in March 2004, was conducted by Earth Tech using the National Functional Guidelines for Organic Data Review, (US EPA, October, 1999), as appropriate. The results were reported by GCAL under SDG 204031909.

| GCAL # | Sample Description |
|-------------|--------------------|
| 20403190901 | SKGW581009 |
| 20403190902 | SKGW58DUP1009 |
| 20403190903 | SKGW631009 |
| 20403190904 | SKGW611009 |
| 20403190905 | SKGW61MS1009 |
| 20403190906 | SKGW61MSD1009 |
| 20403190908 | SKGWFB1009 |
| 20403190909 | SKTB1009 |
| 20403220801 | SKGW06R1009 |
| 20403220802 | SKGW07R1009 |
| 20403220803 | SKGW591009 |
| 20403220804 | SKGW601009 |
| 20403220805 | SKGW62A1009 |
| 20403220806 | SKGW641009 |
| 20403220807 | SKGW651009 |
| 20403220814 | TRIP BLANK |

INTRODUCTION

Analyses were performed according to CLP-Organic Analysis Low Concentration OLC02.0 SOW. Results of the sample analyses are reported by the laboratory as either qualified or unqualified. Unqualified results mean that the reported values may be used without reservation. Various qualifier codes are used by the laboratory to denote specific information regarding the analytical results. The data validation process is intended to evaluate the data on a technical basis. The data package also was subjected to an internal laboratory quality review prior to submission to Earth Tech for data validation.

During the validation process, laboratory-qualified and unqualified data are verified against all available supporting documentation. Based on this evaluation, qualifier codes may be added, deleted or modified by the data user. Final results are therefore, either qualified or unqualified. Validator-qualified results are annotated with the following codes in accordance with the Functional Guidelines:

- U The constituent was analyzed for, but was not detected above the level of the associated analytical reporting limit. The associated value is either the sample quantitation limit or the sample detection limit.
- J The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample.
- UJ The analyte was not detected above the reported sample quantitation limit.
- However, the reported quantitation limit is approximate and may or may not represent the actual limit of quantitation necessary to accurately and precisely measure the analyte in the sample.
- R The sample results are rejected due to serious deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte cannot be verified.

The volatiles data validation findings and conclusions are provided in the following sections of this report:

1. Holding Times
2. GC/MS Tuning
3. Calibration
 - A. IC
 - B. CC
4. Blanks
5. System Monitoring Compound Recovery
6. MS/MSD
7. Laboratory Control Sample
8. Internal Standards Performance
9. Compound Identification
10. Constituent Quantitation and Reported Detection Limits
11. System Performance

- 12. Documentation
- 13. Overall Assessment

1. HOLDING TIMES

All samples for Volatile Organic Compounds (VOC) analyses were analyzed within the 14-day technical holding time and the 10-day VTSR method holding time. The cooler temperature upon receipt at the laboratory was within the recommended temperature of 4°C +/- 2°C.

2. GC/MS TUNING

The samples were analyzed on two GC/MS system, identified as MSV0 and MSV2. Two bromofluorobenzene (BFB) tunes were run. The BFB tunes are acceptable.

3. CALIBRATION

A. Initial Calibration

Two IC's dated 3/21/04 and 3/23/04 were analyzed in support of the volatile sample analyses reported in the data submissions. The IC dated 3/21/04 was analyzed on instrument MSV0 and the IC dated 3/23/04 was analyzed on instrument MSV2. Documentation of the IC standards is present in the data package, and RRF's as well as %RSD values were accurately reported. The criteria employed for technical data review purposes are different than those used in the method. The laboratory must meet a minimum RRF of 0.01; however, for data review purposes, a RRF criterion of "greater than or equal to 0.05" is applied to all volatile compounds. The %RSD's were within the acceptance criteria specified in the method for all target analytes.

The RRF's and the average RRF for the IC dated 3/21/04 were within the acceptance criteria specified in the method for all target analytes with the exception of Acetone and 2-Butnaone. The RRF for the 1.0 ppb standard was below the acceptance criteria. The data validator dropped the 1.0 ppb standard for 1,2-Dibromo-3-chloropropane from the calibration curve and re-calculated the average RRF and %RSD. 1,2-Dibromo-3-chloropropane was not detected in any of the associated samples therefore data qualification was not required.

The RRF's and the average RRF for the IC dated 3/23/04 were within the acceptance criteria specified in the method for all target analytes with the exception of Acetone.

As per the National Functional Guidelines, if any initial calibration RRF is less than 0.05, qualify positive results that have acceptable mass spectral identification with "J", using professional judgement, and non-detected analytes as unusable (R).

It should be noted that the laboratory did meet the minimum RRF of 0.01 for all target compounds.

B. Continuing Calibration

Two CC dated 3/21/04 and 3/23/04 were analyzed on instrument MSV2 in support of the volatile sample analyses reported in the data submissions.

CC dated 3/21/04

The percent difference (%D) between the average RRF's and the CC RF's were within the acceptance criteria for all target analytes. The CC RRF's were within the acceptance criteria specified in the method for all target analytes with the exception of Acetone and 2-Butnaone. The Acetone and 2-Butnaone results were previously qualified under section 3A above.

CC dated 3/23/04

The percent difference (%D) between the average RRF's and the CC RF's were within the acceptance criteria for all target analytes. The CC RRF's were within the acceptance criteria specified in the method for all target analytes with the exception of Acetone. The Acetone results were previously qualified under section 3A above.

4. BLANKS

Two laboratory volatile method blanks, storage blank, two Trip Blanks, and a Field Blank were analyzed with this SDG. The results are summarized below.

MB153850

1,2-Dichlorobenzene and 1,3-Dichlorobenzene were detected at concentrations of 0.20 ppb and 0.12 ppb respectively in the method blank analyzed on 3/21/04.

MB153988

Acetone, 1,2-Dichlorobenzene and 1,3-Dichlorobenzene were detected at concentrations of 1.5 ppb, 0.21 ppb and 0.16 ppb respectively in the method blank analyzed on 3/23/04.

Storage Blank (CLP Storage Blank)

Methylene chloride, 1,4-Dichlorobenzene, and 1,2-Dichlorobenzene were detected at concentrations of 0.15 ppb, 0.096 ppb and 0.11 ppb respectively in the storage blank analyzed on 3/23/04.

Trip Blank (SKTB1009)

Methylene chloride was detected at a concentration of 0.23 ppb in the Trip Blank submitted for the sampling event that occurred on 3/18/04. The Methylene chloride detected in the trip blank was mitigated by the presence of Methylene chloride in the associated storage blank.

Trip Blank

Acetone and Methylene chloride were detected at concentrations 3.1 ppb and 0.24 ppb respectively in the Trip Blank submitted for the sampling event that occurred on 3/16-17/04. The Methylene chloride detected in the trip blank was mitigated by the presence of Methylene chloride in the associated storage blank.

Field Blank (SKGWFB1009)

Methylene chloride was detected at a concentration of 0.52 ppb in the Field Blank collected on 3/18/04. The Methylene chloride detected in the trip blank was mitigated by the presence of Methylene chloride in the associated storage blank.

5. SYSTEM MONITORING COMPOUND RECOVERY

All reported volatile system monitoring compounds were recovered within acceptable control limits for all samples.

6. MATRIX SPIKE/MATRIX SPIKE DUPLICATE

Sample SKGWD611009 was submitted for MS MSD analysis. The MS/MSD percent recoveries (60%-140%) and %RPD (<40%) between the MS MSD were within the acceptance criteria with the exception of the following:

| Compound | MS | MSD | RPD |
|----------------------|-----|-----|-----|
| 1,1-Dichloroethene | 142 | 150 | 5 |
| 2-Hexanone | 58 | 87 | 40 |
| 4-methyl-2-pentanone | 67 | 102 | 41 |
| Bromomethane | 151 | 156 | 3 |
| Carbon disulfide | 145 | 145 | 0 |
| Chloroethane | 132 | 145 | 9 |
| Chloromethane | 135 | 146 | 8 |
| Vinyl chloride | 135 | 143 | 6 |

The results that are bolded exceeded the acceptance criteria. As per the National Functional Guidelines, no action is taken on MS/MSD data alone.

7. LABORATORY CONTROL SAMPLE

A LCS was analyzed in conjunction with this SDG. Recoveries were within the control limit for all constituents.

8. INTERNAL STANDARDS PERFORMANCE

Internal Standard (IS) areas and retention times were within acceptable limits for the reported volatile sample analyses.

9. COMPOUND IDENTIFICATION

All reported VOCs were correctly identified with supporting chromatograms present in the data package.

10. CONSTITUENT QUANTITATION AND REPORTED DETECTION LIMITS

Constituent quantitations were correctly calculated and reported for VOCs with the exception of Ethylbenzene. The Ethylbenzene standard and detected results were originally quantitated using the incorrect quantitation ion (GCAL used 106 instead of 91). GCAL corrected the mistake and re-submitted the corrected pages that were affected in the laboratory report. The overall effect had no impact in the final result for Ethylbenzene.

11. SYSTEM PERFORMANCE

The analytical system appears to have been working well at the time of these analyses, based on the evaluation of the raw data.

12. DOCUMENTATION

The documentation appeared accurate and in order with the exception of the following: GCAL inadvertently transposed the area counts and retention times for IS Chlorobenzene-d5 and 1,4-Dichlorobenzene-d4. The data validator corrected the mistake by drawing arrows to indicate the correct area counts and retention times for IS Chlorobenzene-d5 and 1,4-Dichlorobenzene-d4.

13. OVERALL ASSESSMENT

The results are acceptable with the validator-added qualifiers.

**DATA VALIDATION SUMMARY - SAMPLE DELIVERY GROUP 204031909
PESTICIDES**

Validation of the Gas Chromatography (GC) pesticides data, as prepared by Gulf Coast Analytical Laboratories (GCAL) for the samples collected from the Skinner Landfill site in March 2004, was conducted by Earth Tech using the National Functional Guidelines for Organic Data Review, (US EPA, October, 1999), as appropriate. The results were reported by GCAL under SDG 204031909.

| GCAL # | Sample Description |
|---------------|---------------------------|
| 20403190901 | SKGW581009 |
| 20403190902 | SKGW58DUP1009 |
| 20403190903 | SKGW631009 |
| 20403190904 | SKGW611009 |
| 20403190905 | SKGW61MS1009 |
| 20403190906 | SKGW61MSD1009 |
| 20403190908 | SKGWFB1009 |
| 20403220801 | SKGW06R1009 |
| 20403220802 | SKGW07R1009 |
| 20403220803 | SKGW591009 |
| 20403220804 | SKGW601009 |
| 20403220805 | SKGW62A1009 |
| 20403220806 | SKGW641009 |

INTRODUCTION

Analyses were performed according to CLP-Organic Analysis Multi-Media, Multi-Concentration OLM04.2 SOW. Results of the sample analyses are reported by the laboratory as either qualified or unqualified. Unqualified results mean that the reported values may be used without reservation. Various qualifier codes are used by the laboratory to denote specific information regarding the analytical results.

The data validation process is intended to evaluate the data on a technical basis. The data package also was subjected to an internal laboratory quality review prior to submission to Earth Tech for data validation.

During the validation process, laboratory-qualified and unqualified data are verified against all available supporting documentation. Based on this evaluation, qualifier codes may be added, deleted or modified by the data user. Final results are therefore, either qualified or unqualified. Validator-qualified results are annotated with the following codes in accordance with the Functional Guidelines:

- U** The constituent was analyzed for, but was not detected above the level of the associated analytical reporting limit. The associated value is either the sample quantitation limit or the sample detection limit.

- J The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample.
- UJ The analyte was not detected above the reported sample quantitation limit. However, the reported quantitation limit is approximate and may or may not represent the actual limit of quantitation necessary to accurately and precisely measure the analyte in the sample.
- R The sample results are rejected due to serious deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte cannot be verified.

Details of the pesticide data validation findings and conclusions are provided in the following sections of this report:

1. Holding Times
2. Gas Chromatograph/Electronic Capture Detector (GC/ECD) Instrument Performance Check
3. IC
4. Calibration Verification
5. Blanks
6. Surrogate Spikes
7. Matrix Spike/Matrix Spike Duplicate (MS/MSD)
8. Pesticide Cleanup Checks
9. Target Compound Identification
10. Constituent Quantitation and Reported Detection Limits
11. Documentation
12. Overall Assessment

1. HOLDING TIMES

All samples were extracted within the seven-day ~~technical~~ holding time and the five-day VTSR method holding time. The cooler temperature upon receipt at the laboratory was within the recommended temperature of 4°C +/- 2°C.

2. GC/ECD INSTRUMENT PERFORMANCE CHECK

The Performance Evaluation Mixture (PEM) was analyzed at the correct frequency. Absolute retention times were within limits.

The percent resolution between adjacent peaks was within QC limits for the Pesticide Analyte Resolution Check. The percent resolution between adjacent peaks is within QC limits for the Performance Evaluation Mixtures (PEM). The percent breakdown for both 4,4-DDT and Endrin in each PEM was less than 20.0% for both GC columns. The combined percent breakdown for 4,4-DDT and Endrin in each PEM was less than 30.0% for both GC columns.

3. INITIAL CALIBRATION

Individual standard mixtures A and B were analyzed at the correct frequencies and concentrations. The percent resolution criterion was met for Individual standard mixtures A and B.

The Percent Relative Standard Deviation (%RSD) of the calibration factors for each of the single component pesticides was less than 20%.

The multi-component target compounds were analyzed separately on both columns at a single concentration level. Retention times were determined from a minimum of three peaks.

4. CALIBRATION VERIFICATION

Absolute retention times were within appropriate time retention windows

5. BLANKS

One laboratory method blank and field blank were analyzed with this SDG. The results are summarized below.

Method Blank 154072

No constituents were detected above the laboratory-reporting limit. This blank corresponds to all samples extracted on 3/23/04

Field Blank (SKGWFB1009)

No constituents were detected above the laboratory-reporting limit in the field blank collected on 3/18/04.

6. SURROGATE SPIKES

Decachlorobiphenyl (DCB) and tetrachloro-m-xylene (TCX) surrogate spike recoveries were within the acceptance criteria for all samples.

7. MATRIX SPIKE/MATRIX SPIKE DUPLICATE

Sample SKGWD611009 was submitted for MS/MSD analysis. The MS/MSD percent recoveries were within the acceptance criteria. The %RPD between the MS/MSD are within the acceptance criteria.

8. PESTICIDE CLEANUP CHECKS

Recoveries of all pesticides and surrogates were within 80-120% for the lot of Florisil cartridges utilized for pesticide cleanup with the exception of 4,4'-DDT (130%). There were no target compounds detected in the associated samples therefore no action was taken.

9. TARGET COMPOUND IDENTIFICATION

All reported pesticide data were correctly identified with supporting chromatograms present in the data package.

10. CONSTITUENT QUANTITATION AND REPORTED DETECTION LIMITS

Constituent quantitations were correctly calculated and reported for pesticide constituents.

11. DOCUMENTATION

The documentation appeared accurate and in order with the exception of the following: GCAL inadvertently left the Date Extracted off of the CLP Form 1's therefore the data validator inserted the Date Extracted on the CLP Form 1's.

12. OVERALL ASSESSMENT

The results are acceptable with the validator-added qualifiers.

REFERENCES

US EPA, 1994. *National Functional Guidelines for Inorganic Data Review.*

US EPA, 1999. *National Functional Guidelines for Organic Data Review.*

Me 3/16/04



ANALYTICAL RESULTS

PERFORMED BY

GULF COAST ANALYTICAL LABORATORIES, INC.

Report Date 04/08/2004

GCAL Report 204031909

Deliver To Earth Tech
200 Vine Street
Wilder, KY 41076
859-442-2300

Attn Pat Higgins

Customer Earth Tech

Project Skinner Landfill

000001

CASE NARRATIVE

Client: Earth Tech **Report:** 204031909

Gulf Coast Analytical Laboratories received and analyzed the sample(s) listed on the sample cross-reference page of this report. Receipt of the sample(s) is documented by the attached chain of custody. This applies only to the sample(s) listed in this report. No sample integrity or quality control exceptions were identified unless noted below.

VOLATILES MASS SPECTROMETRY

In analytical batch 271073, the MS MSD exhibited sporadic recovery and RPD failures. The LCS/LCSD recoveries were acceptable.

In analytical batch 271108, no MS MSD was analyzed due to insufficient sample volume. All LCS/LCSD recoveries were acceptable.

SEMI-VOLATILES MASS SPECTROMETRY

The MS/MSD recoveries for 4-Nitrophenol were above the upper control limit. All other batch QC was acceptable.

SEMI-VOLATILE GAS CHROMATOGRAPHY

In the Pesticide analysis, the recovery for DDT was above control limits in the Florisil check, however DDT was not detected in the associated samples.

METALS

In the ILM04.1 - CLP Metals analysis for prep batch 271122, the MS recovery was outside the control limits for Selenium. The LCS recovery was within control limits. This indicates the analysis is in control and the sample is in control and the sample is affected by matrix interference. A post-digestion spike was performed on the QC sample for this batch with a recovery of 49%.

In the ILM04.1 - CLP Metals analysis for prep batch 271124, the MS recovery was outside the control limits for Selenium. The LCS recovery was within control limits. This indicates the analysis is in control and the sample is in control and the sample is affected by matrix interference. A post-digestion spike was performed on the QC sample for this batch with a recovery of 92%.

The Sample/Duplicate RPD for Cyanide, Total for prep batch 271113 is not applicable because the sample and or duplicate concentration is less than five times the reporting limit.

000002

Laboratory Endorsement

Sample analysis was performed in accordance with approved methodologies provided by the Environmental Protection Agency or other recognized agencies. The samples and their corresponding extracts will be maintained for a period of 30 days unless otherwise arranged. Following this retention period the samples will be disposed in accordance with GCAL's Standard Operating Procedures.

Common Abbreviations Utilized in this Report

| | |
|--------------|--|
| ND | Indicates the result was Not Detected at the specified RDL |
| DO | Indicates the result was Diluted Out |
| MI | Indicates the result was subject to Matrix Interference |
| TNTC | Indicates the result was Too Numerous To Count |
| SUBC | Indicates the analysis was Sub-Contracted |
| FLD | Indicates the analysis was performed in the Field |
| PQL | Practical Quantitation Limit |
| MDL | Method Detection Limit |
| RDL | Reporting Detection Limit |
| 00:00 | Reported as a time equivalent to 12:00 AM |

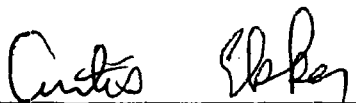
Reporting Flags Utilized in this Report

| | |
|----------|--|
| J | Indicates an estimated value |
| U | Indicates the compound was analyzed for but not detected |
| B | (ORGANICS) Indicates the analyte was detected in the associated Method Blank |
| B | (INORGANICS) Indicates the result is between the RDL and MDL |

Sample receipt at GCAL is documented through the attached chain of custody. In accordance with ISO Guide 25 and NELAC, this report shall be reproduced only in full and with the written permission of GCAL. The results contained within this report relate only to the samples reported. The documented results are presented within this report.

This report pertains only to the samples listed in the Report Sample Summary and should be retained as a permanent record thereof. The results contained within this report are intended for the use of the client. Any unauthorized use of the information contained in this report is prohibited.

I certify that this data package is in compliance with the terms and conditions of the contract and Statement of Work both technically and for completeness, for other than the conditions in the case narrative. Release of the data contained in this hardcopy data package and in the computer-readable data submitted has been authorized by the Quality Assurance Manager or his/her designee, as verified by the following signature.



CURTIS EKKER
DATA VALIDATION MANAGER
GCAL REPORT 204031909

THIS REPORT CONTAINS 627 PAGES.

000003

Report Sample Summary

| GCAL ID | Client ID | Matrix | Collect Date/Time | Receive Date/Time |
|-------------|---------------------|--------|-------------------|-------------------|
| 20403190901 | SKGW581009 | Water | 03/18/2004 11:48 | 03/19/2004 09:45 |
| 20403190902 | SKGW58DUP1009 | Water | 03/18/2004 12:08 | 03/19/2004 09:45 |
| 20403190903 | SKGW631009 | Water | 03/18/2004 14:30 | 03/19/2004 09:45 |
| 20403190904 | SKGW611009 | Water | 03/18/2004 15:35 | 03/19/2004 09:45 |
| 20403190905 | SKGW61MS1009 | Water | 03/18/2004 15:50 | 03/19/2004 09:45 |
| 20403190906 | SKGW61MSD1009 | Water | 03/18/2004 16:10 | 03/19/2004 09:45 |
| 20403190907 | SKGW61DUP1009 | Water | 03/18/2004 16:10 | 03/19/2004 09:45 |
| 20403190908 | SKGWFB1009 | Water | 03/18/2004 17:02 | 03/19/2004 09:45 |
| 20403190909 | SKTB1009 | Water | | 03/19/2004 09:45 |
| 20403190910 | SKGW581009(DISS) | Water | 03/18/2004 11:48 | 03/19/2004 09:45 |
| 20403190911 | SKGW58DUP1009(DISS) | Water | 03/18/2004 12:08 | 03/19/2004 09:45 |
| 20403190912 | SKGW631009(DISS) | Water | 03/18/2004 14:30 | 03/19/2004 09:45 |
| 20403190913 | SKGW611009(DISS) | Water | 03/18/2004 15:35 | 03/19/2004 09:45 |
| 20403190914 | SKGW61MS1009(DISS) | Water | 03/18/2004 15:50 | 03/19/2004 09:45 |
| 20403190915 | SKGW61DUP1009(DISS) | Water | 03/18/2004 16:10 | 03/19/2004 09:45 |
| 20403190916 | SKGWFB1009(DISS) | Water | 03/18/2004 17:02 | 03/19/2004 09:45 |
| 20403220801 | SKGW06R1009 | Water | 03/16/2004 14:58 | 03/19/2004 09:45 |
| 20403220802 | SKGW07R1009 | Water | 03/16/2004 15:52 | 03/19/2004 09:45 |
| 20403220803 | SKGW591009 | Water | 03/17/2004 10:50 | 03/19/2004 09:45 |
| 20403220804 | SKGW601009 | Water | 03/17/2004 11:15 | 03/19/2004 09:45 |
| 20403220805 | SKGW62A1009 | Water | 03/17/2004 11:50 | 03/19/2004 09:45 |
| 20403220806 | SKGW641009 | Water | 03/17/2004 14:10 | 03/19/2004 09:45 |
| 20403220807 | SKGW651009 | Water | 03/17/2004 14:40 | 03/19/2004 09:45 |
| 20403220808 | SKGW06R1009(DISS) | Water | 03/16/2004 14:58 | 03/19/2004 09:45 |
| 20403220809 | SKGW07R1009(DISS) | Water | 03/16/2004 15:52 | 03/19/2004 09:45 |
| 20403220810 | SKGW591009(DISS) | Water | 03/17/2004 10:50 | 03/19/2004 09:45 |
| 20403220811 | SKGW601009(DISS) | Water | 03/17/2004 11:15 | 03/19/2004 09:45 |
| 20403220812 | SKGW62A1009(DISS) | Water | 03/17/2004 11:50 | 03/19/2004 09:45 |
| 20403220813 | SKGW641009(DISS) | Water | 03/17/2004 14:10 | 03/19/2004 09:45 |
| 20403220814 | TRIP BLANK | Water | 03/17/2004 00:00 | 03/19/2004 00:00 |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SKGW581009

Lab Name: GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204031909

Matrix: (soil/water) Water

Sample wt/vol: 25 (g/ml) mL

Lab Sample ID: 20403190901

Level: (low/med) _____

Lab File ID: 2040321/T0692

% Moisture: not dec. _____

Date Collected: 03/18/04 Time: 1148

GC Column: DB-624-30M ID: .53 (mm)

Date Received: 03/19/04

Instrument ID: MSV0

Date Analyzed: 03/21/04 Time: 1941

Soil Extract Volume: _____ (µL)

Dilution Factor: 1 Analyst: RSP

Soil Aliquot Volume: _____ (µL)

Prep Batch: _____ Analytical Batch: 271073

CONCENTRATION UNITS: ug/L

Analytical Method: OLCO 2.1

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|------------|---------------------------|--------|---|-------|-----|
| 71-55-6 | 1,1,1-Trichloroethane | 1.0 | U | 0.010 | 1.0 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1.0 | U | 0.010 | 1.0 |
| 79-00-5 | 1,1,2-Trichloroethane | 1.0 | U | 0.010 | 1.0 |
| 75-34-3 | 1,1-Dichloroethane | 1.0 | U | 0.010 | 1.0 |
| 75-35-4 | 1,1-Dichloroethene | 1.0 | U | 0.010 | 1.0 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 106-93-4 | 1,2-Dibromoethane | 1.0 | U | 0.010 | 1.0 |
| 95-50-1 | 1,2-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 107-06-2 | 1,2-Dichloroethane | 1.0 | U | 0.010 | 1.0 |
| 540-59-0 | 1,2-Dichloroethene | 1.0 | U | 0.010 | 1.0 |
| 78-87-5 | 1,2-Dichloropropane | 1.0 | U | 0.010 | 1.0 |
| 541-73-1 | 1,3-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 106-46-7 | 1,4-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 78-93-3 | 2-Butanone | 5.0 | U | 0.010 | 5.0 |
| 591-78-6 | 2-Hexanone | 5.0 | U | 0.010 | 5.0 |
| 108-10-1 | 4-Methyl-2-pentanone | 5.0 | U | 0.010 | 5.0 |
| 67-64-1 | Acetone | 5.0 | U | 0.010 | 5.0 |
| 71-43-2 | Benzene | 1.0 | U | 0.010 | 1.0 |
| 75-27-4 | Bromodichloromethane | 1.0 | U | 0.010 | 1.0 |
| 75-25-2 | Bromoform | 1.0 | U | 0.010 | 1.0 |
| 74-83-9 | Bromomethane | 1.0 | U | 0.010 | 1.0 |
| 75-15-0 | Carbon disulfide | 1.0 | U | 0.010 | 1.0 |
| 56-23-5 | Carbon tetrachloride | 1.0 | U | 0.010 | 1.0 |
| 108-90-7 | Chlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 75-00-3 | Chloroethane | 1.0 | U | 0.010 | 1.0 |
| 67-66-3 | Chloroform | 1.0 | U | 0.010 | 1.0 |
| 74-87-3 | Chloromethane | 1.0 | U | 0.010 | 1.0 |
| 124-48-1 | Dibromochloromethane | 1.0 | U | 0.010 | 1.0 |
| 10061-01-5 | cis-1,3-Dichloropropene | 1.0 | U | 0.010 | 1.0 |
| 10061-02-6 | trans-1,3-Dichloropropene | 1.0 | U | 0.010 | 1.0 |
| 100-41-4 | Ethylbenzene | 1.0 | U | 0.010 | 1.0 |
| 75-09-2 | Methylene chloride | 2.0 | U | 0.010 | 2.0 |
| 100-42-5 | Styrene | 1.0 | U | 0.010 | 1.0 |

R

R

5/10/04
msc

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SKGW581009

Lab Name: GCAL

Contract:

Lab Code: LA024

Case No.:

SAS No.:

SDG No.: 204031909

Matrix: (soil/water) Water

Sample wt/vol: 25 (g/ml) mL

Lab Sample ID: 20403190901

Level: (low/med)

Lab File ID: 2040321/T0692

% Moisture: not dec.

Date Collected: 03/18/04

Time: 1148

GC Column: DB-624-30M ID: 53 (mm)

Date Received: 03/19/04

Instrument ID: MSV0

Date Analyzed: 03/21/04

Time: 1941

Soil Extract Volume: (µL)

Dilution Factor: 1

Analyst: RSP

Soil Aliquot Volume: (µL)

Prep Batch:

Analytical Batch: 271073

CONCENTRATION UNITS: µg/L

Analytical Method: OLCO 2.1

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|-----------|-------------------|--------|---|-------|-----|
| 127-18-4 | Tetrachloroethene | ND | U | 0.010 | 1.0 |
| 108-88-3 | Toluene | ND | U | 0.010 | 1.0 |
| 79-01-6 | Trichloroethene | ND | U | 0.010 | 1.0 |
| 75-01-4 | Vinyl chloride | ND | U | 0.010 | 1.0 |
| 1330-20-7 | Xylene (total) | ND | U | 0.010 | 1.0 |

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

SKGW581009

Lab Name: GCAL Contract: _____
 Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204031909
 Matrix: Water Lab Sample ID: 20403190901
 Sample wt/vol: _____ Units: _____ Lab File ID: 2040321/T0692
 Level: (low/med) _____ Date Collected: 03/18/04 Time: 1148
 % Moisture: not dec. _____ Date Received: 03/19/04
 GC Column: DB-624-30M ID: .25 (mm) Date Analyzed: 03/21/04 Time: 1941
 Instrument ID: MSV0 Dilution Factor: 1 Analyst: RSP
 Soil Extract Volume: _____ (µL)
 Soil Aliquot Volume: _____ (µL)

Number TICs Found: 0

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | RT | EST. CONC. | Q |
|---------|------------------|----|------------|---|
| 1. | No tics detected | | | |

VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SKGW58DUP1009

Lab Name: GCAL

Contract

Lab Code: LA024

Case No.:

S&S No.

SDG No.: 204031909

Matrix: (soil/water) Water

Sample wt/vol: 25 (g/m) mL

Lab Sample ID: 20403190902

Level: (low/med)

Lab File ID: 2040321/T0688

% Moisture: not dec.

Date Collected: 03/18/04 Time: 1208

GC Column: DB-624-30M ID: 53 (mm)

Date Received: 03/19/04

Instrument ID: MSV0

Date Analyzed: 03/21/04 Time: 1805

Soil Extract Volume: (µL)

Dilution Factor: 1 Analyst: RJO

Soil Aliquot Volume: (µL)

Pres Batch: Analytical Batch: 271073

CONCENTRATION UNITS: µg/L

Analytical Method: OLCO 2.1

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|------------|---------------------------|--------|---|-------|-----|
| 71-55-6 | 1,1,1-Trichloroethane | 1.0 | U | 0.010 | 1.0 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1.0 | U | 0.010 | 1.0 |
| 79-00-5 | 1,1,2-Trichloroethane | 1.0 | U | 0.010 | 1.0 |
| 75-34-3 | 1,1-Dichloroethane | 1.0 | U | 0.010 | 1.0 |
| 75-35-4 | 1,1-Dichloroethene | 1.0 | U | 0.010 | 1.0 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 106-93-4 | 1,2-Dibromoethane | 1.0 | U | 0.010 | 1.0 |
| 95-50-1 | 1,2-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 107-06-2 | 1,2-Dichloroethane | 1.0 | U | 0.010 | 1.0 |
| 540-59-0 | 1,2-Dichloroethene | 1.0 | U | 0.010 | 1.0 |
| 78-87-5 | 1,2-Dichloropropane | 1.0 | U | 0.010 | 1.0 |
| 541-73-1 | 1,3-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 106-46-7 | 1,4-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 78-83-3 | 2-Butanone | 5.0 | U | 0.010 | 5.0 |
| 591-78-6 | 2-Hexanone | 5.0 | U | 0.010 | 5.0 |
| 106-10-1 | 4-Methyl-2-pentanone | 5.0 | U | 0.010 | 5.0 |
| 67-64-1 | Acetone | 5.0 | U | 0.010 | 5.0 |
| 71-43-2 | Benzene | 1.0 | U | 0.010 | 1.0 |
| 75-27-4 | Bromodichloromethane | 1.0 | U | 0.010 | 1.0 |
| 75-25-2 | Bromoform | 1.0 | U | 0.010 | 1.0 |
| 74-83-9 | Bromomethane | 1.0 | U | 0.010 | 1.0 |
| 75-15-0 | Carbon disulfide | 1.0 | U | 0.010 | 1.0 |
| 56-23-5 | Carbon tetrachloride | 1.0 | U | 0.010 | 1.0 |
| 108-90-7 | Chlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 75-00-3 | Chloroethane | 1.0 | U | 0.010 | 1.0 |
| 67-66-3 | Chloroform | 1.0 | U | 0.010 | 1.0 |
| 74-87-3 | Chloromethane | 1.0 | U | 0.010 | 1.0 |
| 124-48-1 | Dibromochloromethane | 1.0 | U | 0.010 | 1.0 |
| 10061-01-5 | cis-1,3-Dichloropropene | 1.0 | U | 0.010 | 1.0 |
| 10061-02-6 | trans-1,3-Dichloropropene | 1.0 | U | 0.010 | 1.0 |
| 100-41-4 | Ethylbenzene | 1.0 | U | 0.010 | 1.0 |
| 75-09-2 | Methylene chloride | 2.0 | U | 0.010 | 2.0 |
| 100-42-5 | Styrene | 1.0 | U | 0.010 | 1.0 |

FORM VOA

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5/21/04
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1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SKGW58DUP1009

Lab Name: GCAL Contract: _____

Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204031909

Matrix: (soil/water) Water

Sample wt/vol: 25 (g/ml) mL

Lab Sample ID: 20403190902

Level: (low/med) _____

Lab File ID: 2040321/T0688

% Moisture: not dec. _____

Date Collected: 03/18/04 Time: 1208

GC Column: DB-624-30M ID: .53 (mm)

Date Received: 03/19/04

Instrument ID: MSV0

Date Analyzed: 03/21/04 Time: 1805

Soil Extract Volume: _____ (μL)

Dilution Factor: 1 Analyst: RJO

Soil Aliquot Volume: _____ (μL)

Prep Batch: _____ Analytical Batch: 271073

CONCENTRATION UNITS: ug/L

Analytical Method: OLCO 2.1

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|-----------|-------------------|--------|---|-------|-----|
| 127-18-4 | Tetrachloroethene | 1.0 | U | 0.010 | 1.0 |
| 108-88-3 | Toluene | 1.0 | U | 0.010 | 1.0 |
| 79-01-6 | Trichloroethene | 1.0 | U | 0.010 | 1.0 |
| 75-01-4 | Vinyl chloride | 1.0 | U | 0.010 | 1.0 |
| 1330-20-7 | Xylene (total) | 1.0 | U | 0.010 | 1.0 |

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

SKGW58DUP1009

Lab Name: GCAL Contract: _____

Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204031909

Matrix: Water Lab Sample ID: 20403190902

Sample wt/vol: _____ Units: _____ Lab File ID: 2040321/T0688

Level: (low/med) _____ Date Collected: 03/18/04 Time: 1208

% Moisture: not dec. _____ Date Received: 03/19/04

GC Column: DB-624-30M ID: 25 (mm) Date Analyzed: 03/21/04 Time: 1805

Instrument ID: MSV0 Dilution Factor: 1 Analyst: RJO

Soil Extract Volume: _____ (µL)

Soil Aliquot Volume: _____ (µL)

Number TICs Found: 0

CONCENTRATION UNITS

| CAS NO. | COMPOUND | RT | EST. CONC. | Q |
|---------|------------------|----|------------|---|
| 1. | No tics detected | | | |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SKGW631009

Lab Name: GCAL

Contract: _____

Lab Code: LA024

Case No.: _____

SAS No.: _____

SDG No.: 204031909

Matrix: (soil/water) Water

Sample wt/vol: 25 (g/ml) mL

Lab Sample ID: 20403190903

Level: (low/med) _____

Lab File ID: 2040321/T0689

% Moisture: not dec. _____

Date Collected: 03/18/04 Time: 1430

GC Column: DB-624-30M ID: .53 (mm)

Date Received: 03/19/04

Instrument ID: MSV0

Date Analyzed: 03/21/04 Time: 1829

Soil Extract Volume: _____ (µL)

Dilution Factor: 1 Analyst: RSP

Soil Aliquot Volume: _____ (µL)

Prep Batch: _____ Analytical Batch: 271073

CONCENTRATION UNITS: ug/L

Analytical Method: OLCO 2.1

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|------------|---------------------------|--------|---|-------|-----|
| 71-55-6 | 1,1,1-Trichloroethane | 1.0 | U | 0.010 | 1.0 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1.0 | U | 0.010 | 1.0 |
| 79-00-5 | 1,1,2-Trichloroethane | 1.0 | U | 0.010 | 1.0 |
| 75-34-3 | 1,1-Dichloroethane | 1.0 | U | 0.010 | 1.0 |
| 75-35-4 | 1,1-Dichloroethene | 1.0 | U | 0.010 | 1.0 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 106-93-4 | 1,2-Dibromoethane | 1.0 | U | 0.010 | 1.0 |
| 95-50-1 | 1,2-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 107-06-2 | 1,2-Dichloroethane | 1.0 | U | 0.010 | 1.0 |
| 540-59-0 | 1,2-Dichloroethene | 1.0 | U | 0.010 | 1.0 |
| 78-87-5 | 1,2-Dichloropropane | 1.0 | U | 0.010 | 1.0 |
| 541-73-1 | 1,3-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 106-46-7 | 1,4-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 78-93-3 | 2-Butanone | 5.0 | U | 0.010 | 5.0 |
| 591-78-6 | 2-Hexanone | 5.0 | U | 0.010 | 5.0 |
| 108-10-1 | 4-Methyl-2-pentanone | 5.0 | U | 0.010 | 5.0 |
| 67-64-1 | Acetone | 5.0 | U | 0.010 | 5.0 |
| 71-43-2 | Benzene | 1.0 | U | 0.010 | 1.0 |
| 75-27-4 | Bromodichloromethane | 1.0 | U | 0.010 | 1.0 |
| 75-25-2 | Bromoform | 1.0 | U | 0.010 | 1.0 |
| 74-83-9 | Bromomethane | 1.0 | U | 0.010 | 1.0 |
| 75-15-0 | Carbon disulfide | 1.0 | U | 0.010 | 1.0 |
| 56-23-5 | Carbon tetrachloride | 1.0 | U | 0.010 | 1.0 |
| 108-90-7 | Chlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 75-00-3 | Chloroethane | 1.0 | U | 0.010 | 1.0 |
| 67-66-3 | Chloroform | 1.0 | U | 0.010 | 1.0 |
| 74-87-3 | Chloromethane | 1.0 | U | 0.010 | 1.0 |
| 124-48-1 | Dibromochloromethane | 1.0 | U | 0.010 | 1.0 |
| 10061-01-5 | cis-1,3-Dichloropropene | 1.0 | U | 0.010 | 1.0 |
| 10061-02-6 | trans-1,3-Dichloropropene | 1.0 | U | 0.010 | 1.0 |
| 100-41-4 | Ethylbenzene | 1.0 | U | 0.010 | 1.0 |
| 75-09-2 | Methylene chloride | 2.0 | U | 0.010 | 2.0 |
| 100-42-5 | Styrene | 1.0 | U | 0.010 | 1.0 |

VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SKGW631009

Lab Name: GCAL Contract: _____

Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204031909

Matrix: (soil/water) Water

Sample wt/vol: 25 (g/ml) mL Lab Sample ID: 20403190903

Level: (low/med) _____ Lab File ID: 2040321/T0689

% Moisture: not dec. _____ Date Collected: 03/18/04 Time: 1430

GC Column: DB-624-30M ID: 53 (mm) Date Received: 03/19/04

Instrument ID: MSV0 Date Analyzed: 03/21/04 Time: 1829

Soil Extract Volume: _____ (µL) Dilution Factor: 1 Analyst: RSP

Soil Aliquot Volume: _____ (µL) Prep Batch: _____ Analytical Batch: 271073

Concentration Units: ug/L Analytical Method: OLCO 2.1

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|-----------|-------------------|--------|---|-------|-----|
| 127-18-4 | Tetrachloroethene | 1.0 | U | 0.010 | 1.0 |
| 108-88-3 | Toluene | 1.0 | U | 0.010 | 1.0 |
| 79-01-6 | Trichloroethene | 1.0 | U | 0.010 | 1.0 |
| 75-01-4 | Vinyl chloride | 1.0 | U | 0.010 | 1.0 |
| 1330-20-7 | Xylene (total) | 1.0 | U | 0.010 | 1.0 |

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

SKGW631009

Lab Name: GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204031909
Matrix: Water Lab Sample ID: 20403190903
Sample wt/vol: _____ Units: _____ Lab File ID: 2040321/T0689
Level: (low/med) _____ Date Collected: 03/18/04 Time: 1430
% Moisture: not dec. _____ Date Received: 03/19/04
GC Column: DB-624-30M ID: .25 (mm) Date Analyzed: 03/21/04 Time: 1829
Instrument ID: MSV0 Dilution Factor: 1 Analyst: RSP
Soil Extract Volume: _____ (μ L)
Soil Aliquot Volume: _____ (μ L)

Number TICs Found: 0

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | RT | EST. CONC. | Q |
|---------|------------------|----|------------|---|
| 1. | No tics detected | | | |

VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SKGW611009

Lab Name: GCAL

Contract: _____

Lab Code: LA024

Case No.: _____

SAS No.: _____

SDG No.: 204031909

Matrix (soil/water) Water

Sample wt/vol: 25 (g/ml) mL

Lab Sample ID: 20403190904

Level: (low/med) _____

Lab File ID: 2040321/T0682

% Moisture: not dec. _____

Date Collected: 03/18/04 Time: 1535

GC Column: DB-624-30M ID: 53 (mm)

Date Received: 03/19/04

Instrument ID: MSV0

Date Analyzed: 03/21/04 Time: 1507

Soil Extract Volume: _____ (µL)

Dilution Factor: 1 Analyst: RJO

Soil Aliquot Volume: _____ (µL)

Prep Batch: _____ Analytical Batch: 271073

CONCENTRATION UNITS: µg/L

Analytical Method: OLCO 2.1

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|------------|---------------------------|--------|---|-------|-----|
| 71-55-6 | 1,1,1-Trichloroethane | 1.0 | U | 0.010 | 1.0 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1.0 | U | 0.010 | 1.0 |
| 79-00-5 | 1,1,2-Trichloroethane | 1.0 | U | 0.010 | 1.0 |
| 75-34-3 | 1,1-Dichloroethane | 1.0 | U | 0.010 | 1.0 |
| 75-35-4 | 1,1-Dichloroethene | 1.0 | U | 0.010 | 1.0 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 106-03-4 | 1,2-Dibromoethane | 1.0 | U | 0.010 | 1.0 |
| 95-50-1 | 1,2-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 107-06-2 | 1,2-Dichloroethane | 1.0 | U | 0.010 | 1.0 |
| 540-59-0 | 1,2-Dichloroethene | 1.0 | U | 0.010 | 1.0 |
| 78-87-5 | 1,2-Dichloropropane | 1.0 | U | 0.010 | 1.0 |
| 541-73-1 | 1,3-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 108-46-7 | 1,4-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 78-83-3 | 2-Butanone | 5.0 | U | 0.010 | 5.0 |
| 591-78-6 | 2-Hexanone | 5.0 | U | 0.010 | 5.0 |
| 108-10-1 | 4-Methyl-2-pentanone | 5.0 | U | 0.010 | 5.0 |
| 67-64-1 | Acetone | 5.0 | U | 0.010 | 5.0 |
| 71-43-2 | Benzene | 1.0 | U | 0.010 | 1.0 |
| 75-27-4 | Bromodichloromethane | 1.0 | U | 0.010 | 1.0 |
| 75-25-2 | Bromoform | 1.0 | U | 0.010 | 1.0 |
| 74-83-9 | Bromomethane | 1.0 | U | 0.010 | 1.0 |
| 75-15-0 | Carbon disulfide | 1.0 | U | 0.010 | 1.0 |
| 56-23-5 | Carbon tetrachloride | 1.0 | U | 0.010 | 1.0 |
| 108-90-7 | Chlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 75-00-3 | Chloroethane | 1.0 | U | 0.010 | 1.0 |
| 67-66-3 | Chloroform | 1.0 | U | 0.010 | 1.0 |
| 74-87-3 | Chloromethane | 1.0 | U | 0.010 | 1.0 |
| 124-48-1 | Dibromochloromethane | 1.0 | U | 0.010 | 1.0 |
| 10061-01-5 | cis-1,3-Dichloropropene | 1.0 | U | 0.010 | 1.0 |
| 10061-02-6 | trans-1,3-Dichloropropene | 1.0 | U | 0.010 | 1.0 |
| 100-41-4 | Ethylbenzene | 1.0 | U | 0.010 | 1.0 |
| 75-09-2 | Methylene chloride | 2.0 | U | 0.010 | 2.0 |
| 100-42-5 | Styrene | 1.0 | U | 0.010 | 1.0 |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SKGW611009

Lab Name: GCAL Contract: _____

Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204031909

Matrix: (soil/water) Water

Sample wt/vol: 25 (g/ml) mL Lab Sample ID: 20403190904

Level: (low/med) _____ Lab File ID: 2040321/T0682

% Moisture: not dec. _____ Date Collected: 03/18/04 Time: 1535

GC Column: DB-624-30M ID: .53 (mm) Date Received: 03/19/04

Instrument ID: MSV0 Date Analyzed: 03/21/04 Time: 1507

Soil Extract Volume: _____ (μ L) Dilution Factor: 1 Analyst: RJO

Soil Aliquot Volume: _____ (μ L) Prep Batch: _____ Analytical Batch: 271073

CONCENTRATION UNITS: ug/L Analytical Method: OLCO 2.1

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|-----------|-------------------|--------|---|-------|-----|
| 127-18-4 | Tetrachloroethene | 1.0 | U | 0.010 | 1.0 |
| 108-88-3 | Toluene | 1.0 | U | 0.010 | 1.0 |
| 79-01-6 | Trichloroethene | 1.0 | U | 0.010 | 1.0 |
| 75-01-4 | Vinyl chloride | 1.0 | U | 0.010 | 1.0 |
| 1330-20-7 | Xylene (total) | 1.0 | U | 0.010 | 1.0 |

VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

SKGW611009

Lab Name: GCAL Contract: _____

Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204031909

Matrix: Water Lab Sample ID: 20403190904

Sample wt/vol: _____ Units: _____ Lab File ID: 2040321/T0682

Level: (low/med) _____ Date Collected: 03/18/04 Time: 1535

% Moisture: not dec. _____ Date Received: 03/19/04

GC Column: DB-624-30M ID: 25 (mm) Date Analyzed: 03/21/04 Time: 1507

Instrument ID: MSV0 Dilution Factor: 1 Analyst: RSP

Soil Extract Volume: _____ (µL)

Soil Aliquot Volume: _____ (µL)

Number TICs Found: 0

CONCENTRATION UNITS

| CAS NO. | COMPOUND | RT | EST. CONC. | Q |
|---------|------------------|----|------------|---|
| 1. | No tics detected | | | |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SKGWFB1009

Lab Name: GCAL

Contract:

Lab Code: LA024

Case No.:

SAS No.:

SDG No.: 204031909

Matrix: (soil/water) Water

Sample wt/vol: 25 (g/ml) mL

Lab Sample ID: 20403190908

Level: (low/med)

Lab File ID: 2040321/T0690

% Moisture: not dec.

Date Collected: 03/18/04

Time: 1702

GC Column: DB-624-30M

ID: .53 (mm)

Date Received: 03/19/04

Instrument ID: MSV0

Date Analyzed: 03/21/04

Time: 1853

Soil Extract Volume: (µL)

Dilution Factor: 1

Analyst: RSP

Soil Aliquot Volume: (µL)

Prep Batch:

Analytical Batch: 271073

CONCENTRATION UNITS: ug/L

Analytical Method: OLCO 2.1

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|------------|---------------------------|----------|---|-------|-----|
| 71-55-6 | 1,1,1-Trichloroethane | 1.0 | U | 0.010 | 1.0 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1.0 | U | 0.010 | 1.0 |
| 79-00-5 | 1,1,2-Trichloroethane | 1.0 | U | 0.010 | 1.0 |
| 75-34-3 | 1,1-Dichloroethane | 1.0 | U | 0.010 | 1.0 |
| 75-35-4 | 1,1-Dichloroethene | 1.0 | U | 0.010 | 1.0 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 106-93-4 | 1,2-Dibromoethane | 1.0 | U | 0.010 | 1.0 |
| 95-50-1 | 1,2-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 107-06-2 | 1,2-Dichloroethane | 1.0 | U | 0.010 | 1.0 |
| 540-59-0 | 1,2-Dichloroethene | 1.0 | U | 0.010 | 1.0 |
| 78-87-5 | 1,2-Dichloropropane | 1.0 | U | 0.010 | 1.0 |
| 541-73-1 | 1,3-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 106-46-7 | 1,4-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 78-93-3 | 2-Butanone | 5.0 | U | 0.010 | 5.0 |
| 591-78-6 | 2-Hexanone | 5.0 | U | 0.010 | 5.0 |
| 108-10-1 | 4-Methyl-2-pentanone | 5.0 | U | 0.010 | 5.0 |
| 67-64-1 | Acetone | 5.0 | U | 0.010 | 5.0 |
| 71-43-2 | Benzene | 1.0 | U | 0.010 | 1.0 |
| 75-27-4 | Bromodichloromethane | 1.0 | U | 0.010 | 1.0 |
| 75-25-2 | Bromoform | 1.0 | U | 0.010 | 1.0 |
| 74-83-9 | Bromomethane | 1.0 | U | 0.010 | 1.0 |
| 75-15-0 | Carbon disulfide | 1.0 | U | 0.010 | 1.0 |
| 56-23-5 | Carbon tetrachloride | 1.0 | U | 0.010 | 1.0 |
| 108-90-7 | Chlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 75-00-3 | Chloroethane | 1.0 | U | 0.010 | 1.0 |
| 67-66-3 | Chloroform | 1.0 | U | 0.010 | 1.0 |
| 74-87-3 | Chloromethane | 1.0 | U | 0.010 | 1.0 |
| 124-48-1 | Dibromochloromethane | 1.0 | U | 0.010 | 1.0 |
| 10061-01-5 | cis-1,3-Dichloropropene | 1.0 | U | 0.010 | 1.0 |
| 10061-02-6 | trans-1,3-Dichloropropene | 1.0 | U | 0.010 | 1.0 |
| 100-41-4 | Ethylbenzene | 1.0 | U | 0.010 | 1.0 |
| 75-09-2 | Methylene chloride | 2.0 0.52 | J | 0.010 | 2.0 |
| 100-42-5 | Styrene | 1.0 | U | 0.010 | 1.0 |

*A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SKGWFB1009

Lab Name: GCAL Contract: _____

Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204031909

Matrix (soil/water) Water

Sample wt/vol: 25 (g/ml) mL Lab Sample ID: 20403190908

Level: (low/med) _____ Lab File ID: 2040321/T0690

% Moisture: not dec. _____ Date Collected: 03/18/04 Time: 1702

GC Column: DB-624-30M ID: 53 (mm) Date Received: 03/19/04

Instrument ID: MSV0 Date Analyzed: 03/21/04 Time: 1853

Soil Extract Volume: _____ (µL) Dilution Factor: 1 Analyst: RSP

Soil Aliquot Volume: _____ (µL) Prep Batch: _____ Analytical Batch: 271073

Analytical Method: OLCO 2.1

CONCENTRATION UNITS µg/L

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|-----------|-------------------|--------|---|-------|-----|
| 127-18-4 | Tetrachloroethene | 1.0 | U | 0.010 | 1.0 |
| 108-88-3 | Toluene | 1.0 | U | 0.010 | 1.0 |
| 79-01-6 | Trichloroethene | 1.0 | U | 0.010 | 1.0 |
| 75-01-4 | Vinyl chloride | 1.0 | U | 0.010 | 1.0 |
| 1330-20-7 | Xylene (total) | 1.0 | U | 0.010 | 1.0 |

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

SKGWFB1009

Lab Name: GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204031909
Matrix: Water Lab Sample ID: 20403190908
Sample wt/vol: _____ Units: _____ Lab File ID: 2040321/T0690
Level: (low/med) _____ Date Collected: 03/18/04 Time: 1702
% Moisture: not dec. _____ Date Received: 03/19/04
GC Column: DB-624-30M ID: .53 (mm) Date Analyzed: 03/21/04 Time: 1853
Instrument ID: MSV0 Dilution Factor: 1 Analyst: RSP
Soil Extract Volume: _____ (µL)
Soil Aliquot Volume: _____ (µL)

Number TICs Found: 1

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | RT | EST. CONC. | Q |
|--------------|----------------|-------|------------|---|
| 1. 7446-09-5 | Sulfur dioxide | 1.953 | 197 | |

*A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SKTB1009

Lab Name: GCAL

Contract: _____

Lab Code: LA024

Case No.: _____

SAS No.: _____

SDG No.: 204031909

Matrix: (soil/water) Water

Sample wt/vol: 25 (g/ml) ml

Lab Sample ID: 20403190909

Level: (low/med) _____

Lab File ID: 2040321/T0691

% Moisture: not dec. _____

Date Collected: _____ Time: _____

GC Column: DB-624-30M ID: .53 (mm)

Date Received: 03/19/04

Instrument ID: MSV0

Date Analyzed: 03/21/04 Time: 1916

Soil Extract Volume: _____ (μL)

Dilution Factor: 1 Analyst: RSP

Soil Aliquot Volume: _____ (μL)

Prep Batch: _____ Analytical Batch: 271073

CONCENTRATION UNITS μg/L

Analytical Method: OLCO 2.1

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|------------|---------------------------|--------|---|-------|-----|
| 71-55-6 | 1,1,1-Trichloroethane | 1.0 | U | 0.010 | 1.0 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1.0 | U | 0.010 | 1.0 |
| 79-00-5 | 1,1,2-Trichloroethane | 1.0 | U | 0.010 | 1.0 |
| 75-34-3 | 1,1-Dichloroethane | 1.0 | U | 0.010 | 1.0 |
| 75-35-4 | 1,1-Dichloroethane | 1.0 | U | 0.010 | 1.0 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 106-93-4 | 1,2-Dibromoethane | 1.0 | U | 0.010 | 1.0 |
| 95-50-1 | 1,2-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 107-06-2 | 1,2-Dichloroethane | 1.0 | U | 0.010 | 1.0 |
| 540-59-0 | 1,2-Dichloroethane | 1.0 | U | 0.010 | 1.0 |
| 78-87-5 | 1,2-Dichloropropane | 1.0 | U | 0.010 | 1.0 |
| 541-73-1 | 1,3-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 108-46-7 | 1,4-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 78-83-3 | 2-Butanone | 5.0 | U | 0.010 | 5.0 |
| 591-78-6 | 2-Hexanone | 5.0 | U | 0.010 | 5.0 |
| 105-10-1 | 4-Methyl-2-pentanone | 5.0 | U | 0.010 | 5.0 |
| 67-64-1 | Acetone | 5.0 | U | 0.010 | 5.0 |
| 71-43-2 | Benzene | 1.0 | U | 0.010 | 1.0 |
| 75-27-4 | Bromodichloromethane | 1.0 | U | 0.010 | 1.0 |
| 75-25-2 | Bromoform | 1.0 | U | 0.010 | 1.0 |
| 74-83-9 | Bromomethane | 1.0 | U | 0.010 | 1.0 |
| 75-15-0 | Carbon disulfide | 1.0 | U | 0.010 | 1.0 |
| 56-23-5 | Carbon tetrachloride | 1.0 | U | 0.010 | 1.0 |
| 108-90-7 | Chlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 75-00-3 | Chloroethane | 1.0 | U | 0.010 | 1.0 |
| 67-66-3 | Chloroform | 1.0 | U | 0.010 | 1.0 |
| 74-87-3 | Chloromethane | 1.0 | U | 0.010 | 1.0 |
| 124-48-1 | Dibromochloromethane | 1.0 | U | 0.010 | 1.0 |
| 10061-01-5 | cis-1,3-Dichloropropene | 1.0 | U | 0.010 | 1.0 |
| 10061-02-6 | trans-1,3-Dichloropropene | 1.0 | U | 0.010 | 1.0 |
| 100-41-4 | Ethylbenzene | 1.0 | U | 0.010 | 1.0 |
| 75-08-2 | Methylene chloride | 1.0 | J | 0.010 | 2.0 |
| 100-42-5 | Styrene | 1.0 | U | 0.010 | 1.0 |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SKTB1009

Lab Name: GCAL Contract: _____

Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204031909

Matrix: (soil/water) Water

Sample wt/vol: 25 (g/ml) mL Lab Sample ID: 20403190909

Level: (low/med) _____ Lab File ID: 2040321/T0691

% Moisture: not dec. _____ Date Collected: _____ Time: _____

GC Column: DB-624-30M ID: .53 (mm) Date Received: 03/19/04

Instrument ID: MSV0 Date Analyzed: 03/21/04 Time: 1916

Soil Extract Volume: _____ (µL) Dilution Factor: 1 Analyst: RSP

Soil Allquot Volume: _____ (µL) Prep Batch: _____ Analytical Batch: 271073

CONCENTRATION UNITS: ug/L Analytical Method: OLCO 2.1

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|-----------|-------------------|--------|---|-------|-----|
| 127-18-4 | Tetrachloroethene | 1.0 | U | 0.010 | 1.0 |
| 108-88-3 | Toluene | 1.0 | U | 0.010 | 1.0 |
| 79-01-6 | Trichloroethene | 1.0 | U | 0.010 | 1.0 |
| 75-01-4 | Vinyl chloride | 1.0 | U | 0.010 | 1.0 |
| 1330-20-7 | Xylene (total) | 1.0 | U | 0.010 | 1.0 |

VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.
SKTB1009

Lab Name: GCAL Contract: _____
 Lab Code: LA024 Case No.: _____ SAS No. _____ SDG No.: 204031909
 Matrix: Water Lab Sample ID: 20403190909
 Sample wt/vol: _____ Units: _____ Lab File ID: 2040321/T0691
 Level: (low/med) _____ Date Collected: _____ Time: _____
 % Moisture: not dec. _____ Date Received: 03/19/04
 GC Column: DB-624-30M ID: .53 (mm) Date Analyzed: 03/21/04 Time: 1916
 Instrument ID: MSV0 Dilution Factor: 1 Analyst: RSP
 Soil Extract Volume: _____ (µL)
 Soil Aliquot Volume: _____ (µL)

Number TICs Found: _____

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | RT | EST. CONC. | Q |
|--------------|----------------|------|------------|---|
| 1. 7446-09-5 | Sulfur dioxide | 8.78 | 8.77 | |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SKGW06R1009

Lab Name: GCAL

Contract: _____

Lab Code: LA024

Case No.: _____

SAS No.: _____

SDG No.: 204031909

Matrix: (soil/water) Water

Sample wt/vol: 25 (g/ml) mL

Lab Sample ID: 20403220801

Level: (low/med) _____

Lab File ID: 2040323/T2558

% Moisture: not dec. _____

Date Collected: 03/16/04 Time: 1458

GC Column: DB-624-30M ID: .53 (mm)

Date Received: 03/19/04

Instrument ID: MSV2

Date Analyzed: 03/23/04 Time: 1326

Soil Extract Volume: _____ (µL)

Dilution Factor: 1 Analyst: RSP

Soil Aliquot Volume: _____ (µL)

Prep Batch: _____ Analytical Batch: 271108

CONCENTRATION UNITS: ug/L

Analytical Method: OLCO 2.1

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|------------|---------------------------|--------|---|-------|-----|
| 71-55-6 | 1,1,1-Trichloroethane | 1.0 | U | 0.010 | 1.0 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1.0 | U | 0.010 | 1.0 |
| 79-00-5 | 1,1,2-Trichloroethane | 1.0 | U | 0.010 | 1.0 |
| 75-34-3 | 1,1-Dichloroethane | 1.0 | U | 0.010 | 1.0 |
| 75-35-4 | 1,1-Dichloroethene | 1.0 | U | 0.010 | 1.0 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 106-93-4 | 1,2-Dibromoethane | 1.0 | U | 0.010 | 1.0 |
| 95-50-1 | 1,2-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 107-06-2 | 1,2-Dichloroethane | 1.0 | U | 0.010 | 1.0 |
| 540-59-0 | 1,2-Dichloroethene | 1.0 | U | 0.010 | 1.0 |
| 78-87-5 | 1,2-Dichloropropane | 1.0 | U | 0.010 | 1.0 |
| 541-73-1 | 1,3-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 106-46-7 | 1,4-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 78-93-3 | 2-Butanone | 5.0 | U | 0.010 | 5.0 |
| 591-78-6 | 2-Hexanone | 5.0 | U | 0.010 | 5.0 |
| 108-10-1 | 4-Methyl-2-pentanone | 5.0 | U | 0.010 | 5.0 |
| 67-84-1 | Acetone | 5.0 | U | 0.010 | 5.0 |
| 71-43-2 | Benzene | 1.0 | U | 0.010 | 1.0 |
| 75-27-4 | Bromodichloromethane | 1.0 | U | 0.010 | 1.0 |
| 75-25-2 | Bromoform | 1.0 | U | 0.010 | 1.0 |
| 74-83-9 | Bromomethane | 1.0 | U | 0.010 | 1.0 |
| 75-15-0 | Carbon disulfide | 1.0 | U | 0.010 | 1.0 |
| 58-23-5 | Carbon tetrachloride | 1.0 | U | 0.010 | 1.0 |
| 108-90-7 | Chlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 75-00-3 | Chloroethane | 1.0 | U | 0.010 | 1.0 |
| 67-86-3 | Chloroform | 1.0 | U | 0.010 | 1.0 |
| 74-87-3 | Chloromethane | 1.0 | U | 0.010 | 1.0 |
| 124-48-1 | Dibromochloromethane | 1.0 | U | 0.010 | 1.0 |
| 10061-01-5 | cis-1,3-Dichloropropene | 1.0 | U | 0.010 | 1.0 |
| 10061-02-6 | trans-1,3-Dichloropropene | 1.0 | U | 0.010 | 1.0 |
| 100-41-4 | Ethylbenzene | 1.0 | U | 0.010 | 1.0 |
| 75-09-2 | Methylene chloride | 2.0 | J | 0.010 | 2.0 |
| 100-42-5 | Styrene | 1.0 | U | 0.010 | 1.0 |

5/12/04
msw

VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SKGW06R1009

Lab Name: GCAL

Contract: _____

Lab Code: LA024

Case No.: _____

SAS No. _____

SDG No.: 204031909

Matrix: (soil/water) Water

Sample wt/vol: 25 (g/m) mL

Lab Sample ID: 20403220801

Level: (pow/med) _____

Lab File ID: 2040323/T2558

% Moisture: not dec. _____

Date Collected: 03/16/04

Time: 1458

GC Column: DB-624-30M ID: 53 (mm)

Date Received: 03/19/04

Instrument ID: MSV2

Date Analyzed: 03/23/04

Time: 1326

Soil Extract Volume: _____ (µL)

Dilution Factor: 1

Analyst: RSP

Soil Aliquot Volume: _____ (µL)

Prep Batch: _____

Analytical Batch: 271108

CONCENTRATION UNITS: µg/L

Analytical Method: OLCO 2.1

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|-----------|-------------------|--------|---|-------|-----|
| 127-18-4 | Tetrachloroethene | 1.0 | U | 0.010 | 1.0 |
| 108-88-3 | Toluene | 1.0 | U | 0.010 | 1.0 |
| 79-01-6 | Trichloroethene | 1.0 | U | 0.010 | 1.0 |
| 75-01-4 | Vinyl chloride | 1.0 | U | 0.010 | 1.0 |
| 1330-20-7 | Xylene (total) | 1.0 | U | 0.010 | 1.0 |

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

SKGW06R1009

Lab Name: GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204031909
Matrix: Water Lab Sample ID: 20403220801
Sample wt/vol: _____ Units: _____ Lab File ID: 2040321/T2558
Level: (low/med) _____ Date Collected: 03/16/04 Time: 1458
% Moisture: not dec. _____ Date Received: 03/19/04
GC Column: DB-624-30M ID: .25 (mm) Date Analyzed: 03/23/04 Time: 1327
Instrument ID: MSV2 Dilution Factor: 1 Analyst: RSP
Soil Extract Volume: _____ (µL)
Soil Aliquot Volume: _____ (µL)

Number TICs Found: 0

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | RT | EST. CONC. | Q |
|---------|------------------|----|------------|---|
| 1. | No tics detected | | | |

VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SKGW07R1009

Lab Name: GCAL

Contract:

Lab Code: LA024

Case No.:

SAS No.:

SDG No.: 204031909

Matrix: (soil/water) Water

Sample wt/vol: 25 (g/ml) mL

Lab Sample ID: 20403220802

Level: (low/med)

Lab File ID: 2040323/T2559

% Moisture: not dec.

Date Collected: 03/16/04

Time: 1552

GC Column: DB-624-30M

ID: 53

(mm)

Date Received: 03/19/04

Instrument ID: MSV2

Date Analyzed: 03/23/04

Time: 1351

Soil Extract Volume: (µL)

Dilution Factor: 1

Analyst: RSP

Soil Aliquot Volume: (µL)

Prep Batch:

Analytical Batch: 271108

CONCENTRATION UNITS: µg/L

Analytical Method: OLCO 2.1

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|------------|---------------------------|--------|---|-------|-----|
| 71-55-6 | 1,1,1-Trichloroethane | 1.0 | U | 0.010 | 1.0 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1.0 | U | 0.010 | 1.0 |
| 79-00-5 | 1,1,2-Trichloroethane | 1.0 | U | 0.010 | 1.0 |
| 75-34-3 | 1,1-Dichloroethane | 1.0 | U | 0.010 | 1.0 |
| 75-35-4 | 1,1-Dichloroethene | 1.0 | U | 0.010 | 1.0 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 106-83-4 | 1,2-Dibromoethane | 1.0 | U | 0.010 | 1.0 |
| 95-50-1 | 1,2-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 107-06-2 | 1,2-Dichloroethane | 1.0 | U | 0.010 | 1.0 |
| 540-59-0 | 1,2-Dichloroethene | 1.0 | U | 0.010 | 1.0 |
| 78-87-5 | 1,2-Dichloropropane | 1.0 | U | 0.010 | 1.0 |
| 541-73-1 | 1,3-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 106-46-7 | 1,4-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 78-83-3 | 2-Butanone | 5.0 | U | 0.010 | 5.0 |
| 591-78-6 | 2-Hexanone | 5.0 | U | 0.010 | 5.0 |
| 108-10-1 | 4-Methyl-2-pentanone | 5.0 | U | 0.010 | 5.0 |
| 67-64-1 | Acetone | 5.0 | U | 0.010 | 5.0 |
| 71-43-2 | Benzene | 1.0 | U | 0.010 | 1.0 |
| 75-27-4 | Bromodichloromethane | 1.0 | U | 0.010 | 1.0 |
| 75-25-2 | Bromoforn | 1.0 | U | 0.010 | 1.0 |
| 74-83-9 | Bromomethane | 1.0 | U | 0.010 | 1.0 |
| 75-15-0 | Carbon disulfide | 1.0 | U | 0.010 | 1.0 |
| 56-23-5 | Carbon tetrachloride | 1.0 | U | 0.010 | 1.0 |
| 108-90-7 | Chlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 75-00-3 | Chloroethane | 1.0 | U | 0.010 | 1.0 |
| 67-66-3 | Chloroform | 1.0 | U | 0.010 | 1.0 |
| 74-87-3 | Chloromethane | 1.0 | U | 0.010 | 1.0 |
| 124-48-1 | Dibromochloromethane | 1.0 | U | 0.010 | 1.0 |
| 10061-01-5 | cis-1,3-Dichloropropene | 1.0 | U | 0.010 | 1.0 |
| 10061-02-6 | trans-1,3-Dichloropropene | 1.0 | U | 0.010 | 1.0 |
| 100-41-4 | Ethylbenzene | 1.0 | U | 0.010 | 1.0 |
| 75-09-2 | Methylene chloride | 2.0 | J | 0.010 | 2.0 |
| 100-42-5 | Styrene | 1.0 | U | 0.010 | 1.0 |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SKGW07R1009

Lab Name: GCAL Contract: _____

Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204031909

Matrix: (soil/water) Water

Sample wt/vol: 25 (g/ml) mL

Lab Sample ID: 20403220802

Level: (low/med) _____

Lab File ID: 2040323/T2559

% Moisture: not dec. _____

Date Collected: 03/16/04 Time: 1552

GC Column: DB-624-30M ID: .53 (mm)

Date Received: 03/19/04

Instrument ID: MSV2

Date Analyzed: 03/23/04 Time: 1351

Soil Extract Volume: _____ (µL)

Dilution Factor: 1 Analyst: RSP

Soil Aliquot Volume: _____ (µL)

Prep Batch: _____ Analytical Batch: 271108

CONCENTRATION UNITS: ug/L

Analytical Method: OLCO 2.1

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|-----------|-------------------|--------|---|-------|-----|
| 127-18-4 | Tetrachloroethene | 1.0 | U | 0.010 | 1.0 |
| 108-88-3 | Toluene | 1.0 | U | 0.010 | 1.0 |
| 79-01-6 | Trichloroethene | 1.0 | U | 0.010 | 1.0 |
| 75-01-4 | Vinyl chloride | 1.0 | U | 0.010 | 1.0 |
| 1330-20-7 | Xylene (total) | 1.0 | U | 0.010 | 1.0 |

VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

SKGW07R1009

Lab Name: GCAL Contract: _____

Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204031909

Matrix: Water Lab Sample ID: 20403220802

Sample wt/vol: _____ Units: _____ Lab File ID: 2040323/T2559

Level: (low/med) _____ Date Collected: 03/16/04 Time: 1552

% Moisture: not dec. _____ Date Received: 03/19/04

GC Column: DB-624-30M ID: .25 (mm) Date Analyzed: 03/23/04 Time: 1351

Instrument ID: MSV2 Dilution Factor: 1 Analyst: RSP

Soil Extract Volume: _____ (µL)

Soil Aliquot Volume: _____ (µL)

Number TICs Found: 0

CONCENTRATION UNITS

| CAS NO. | COMPOUND | RT | EST. CONC. | Q |
|---------|------------------|----|------------|---|
| 1. | No tics detected | | | |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SKGW591009

Lab Name: GCAL Contract: _____

Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204031909

Matrix: (soil/water) Water

Sample wt/vol: 25 (g/ml) mL

Lab Sample ID: 20403220803

Level: (low/med) _____

Lab File ID: 2040323/T2561

% Moisture: not dec. _____

Date Collected: 03/17/04 Time: 1050

GC Column: DB-624-30M ID: .53 (mm)

Date Received: 03/19/04

Instrument ID: MSV2

Date Analyzed: 03/23/04 Time: 1440

Soil Extract Volume: _____ (µL)

Dilution Factor: 1 Analyst: RSP

Soil Aliquot Volume: _____ (µL)

Prep Batch: _____ Analytical Batch: 271108

CONCENTRATION UNITS: ug/L

Analytical Method: OLCO 2.1

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|------------|---------------------------|--------|---|-------|-----|
| 71-55-6 | 1,1,1-Trichloroethane | 1.0 | U | 0.010 | 1.0 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1.0 | U | 0.010 | 1.0 |
| 79-00-5 | 1,1,2-Trichloroethane | 1.0 | U | 0.010 | 1.0 |
| 75-34-3 | 1,1-Dichloroethane | 1.0 | U | 0.010 | 1.0 |
| 75-35-4 | 1,1-Dichloroethene | 1.0 | U | 0.010 | 1.0 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 106-93-4 | 1,2-Dibromoethane | 1.0 | U | 0.010 | 1.0 |
| 95-50-1 | 1,2-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 107-06-2 | 1,2-Dichloroethane | 1.0 | U | 0.010 | 1.0 |
| 540-59-0 | 1,2-Dichloroethene | 1.0 | U | 0.010 | 1.0 |
| 78-87-5 | 1,2-Dichloropropane | 1.0 | U | 0.010 | 1.0 |
| 541-73-1 | 1,3-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 106-46-7 | 1,4-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 78-93-3 | 2-Butanone | 5.0 | U | 0.010 | 5.0 |
| 591-78-6 | 2-Hexanone | 5.0 | U | 0.010 | 5.0 |
| 108-10-1 | 4-Methyl-2-pentanone | 5.0 | U | 0.010 | 5.0 |
| 67-64-1 | Acetone | 5.0 | U | 0.010 | 5.0 |
| 71-43-2 | Benzene | 1.0 | U | 0.010 | 1.0 |
| 75-27-4 | Bromodichloromethane | 1.0 | U | 0.010 | 1.0 |
| 75-25-2 | Bromoform | 1.0 | U | 0.010 | 1.0 |
| 74-83-9 | Bromomethane | 1.0 | U | 0.010 | 1.0 |
| 75-15-0 | Carbon disulfide | 1.0 | U | 0.010 | 1.0 |
| 56-23-5 | Carbon tetrachloride | 1.0 | U | 0.010 | 1.0 |
| 108-90-7 | Chlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 75-00-3 | Chloroethane | 1.0 | U | 0.010 | 1.0 |
| 67-66-3 | Chloroform | 1.0 | U | 0.010 | 1.0 |
| 74-87-3 | Chloromethane | 1.0 | U | 0.010 | 1.0 |
| 124-48-1 | Dibromochloromethane | 1.0 | U | 0.010 | 1.0 |
| 10061-01-5 | cis-1,3-Dichloropropene | 1.0 | U | 0.010 | 1.0 |
| 10061-02-6 | trans-1,3-Dichloropropene | 1.0 | U | 0.010 | 1.0 |
| 100-41-4 | Ethylbenzene | 1.0 | U | 0.010 | 1.0 |
| 75-09-2 | Methylene chloride | 2.0020 | J | 0.010 | 2.0 |
| 100-42-5 | Styrene | 1.0 | U | 0.010 | 1.0 |

VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SKGW591009

Lab Name: GCAL

Contract: _____

Lab Code: LA024

Case No.: _____

SAS No.: _____

SDG No.: 204031909

Matrix: (soil/water) Water

Sample wt/vol: 25 (g/ml) mL

Lab Sample ID: 20403220803

Level: (low/med) _____

Lab File ID: 2040323/T2561

% Moisture: not dec. _____

Date Collected: 03/17/04 Time: 1050

GC Column: DB-624-30M ID: 53 (mm)

Date Received: 03/19/04

Instrument ID: MSV2

Date Analyzed: 03/23/04 Time: 1440

Soil Extract Volume: _____ (μL)

Dilution Factor: 1 Analyst: RSP

Soil Aliquot Volume: _____ (μL)

Prep Batch: _____ Analytical Batch: 271108

CONCENTRATION UNITS ug/L

Analytical Method: OLCO 2.1

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|-----------|-------------------|--------|---|-------|-----|
| 127-18-4 | Tetrachloroethene | 1.0 | U | 0.010 | 1.0 |
| 108-88-3 | Toluene | 1.0 | U | 0.010 | 1.0 |
| 79-01-8 | Trichloroethene | 1.0 | U | 0.010 | 1.0 |
| 75-01-4 | Vinyl chloride | 1.0 | U | 0.010 | 1.0 |
| 1330-20-7 | Xylene (total) | 1.0 | U | 0.010 | 1.0 |

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

SKGW591009

Lab Name: GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204031909
Matrix: Water Lab Sample ID: 20403220803
Sample wt/vol: _____ Units: _____ Lab File ID: 2040323/T2561
Level: (low/med) _____ Date Collected: 03/17/04 Time: 1050
% Moisture: not dec. _____ Date Received: 03/19/04
GC Column: DB-624-30M ID: .25 (mm) Date Analyzed: 03/23/04 Time: 1440
Instrument ID: MSV2 Dilution Factor: 1 Analyst: RSP
Soil Extract Volume: _____ (µL)
Soil Aliquot Volume: _____ (µL)

Number TICs Found: 0

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | RT | EST. CONC. | Q |
|---------|------------------|----|------------|---|
| 1. | No tics detected | | | |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SKGW601009

Lab Name: GCAL

Contract:

Lab Code: LA024

Case No.:

SAS No.:

SDG No.: 204031909

Matrix (soil/water) Water

Sample wt/vol: 25 (g/ml) mL

Lab Sample ID: 20403220804

Level: (low/med)

Lab File ID: 2040323/T2562

% Moisture: not dec.

Date Collected: 03/17/04

Time: 1115

GC Column: DB-624-30M ID: 53 (mm)

Date Received: 03/19/04

Instrument ID: MSV2

Date Analyzed: 03/23/04

Time: 1505

Soil Extract Volume: (µL)

Dilution Factor: 1

Analyst: RSP

Soil Aliquot Volume: (µL)

Prep Batch:

Analytical Batch: 271108

CONCENTRATION UNITS: µg/L

Analytical Method: OLCO 2.1

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|------------|---------------------------|--------|----|-------|-----|
| 71-55-6 | 1,1,1-Trichloroethane | 1.0 | U | 0.010 | 1.0 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1.0 | U | 0.010 | 1.0 |
| 79-00-5 | 1,1,2-Trichloroethane | 1.0 | U | 0.010 | 1.0 |
| 75-34-3 | 1,1-Dichloroethane | 1.0 | U | 0.010 | 1.0 |
| 75-35-4 | 1,1-Dichloroethene | 1.0 | U | 0.010 | 1.0 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 106-93-4 | 1,2-Dibromoethane | 1.0 | U | 0.010 | 1.0 |
| 95-50-1 | 1,2-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 107-06-2 | 1,2-Dichloroethane | 1.0 | U | 0.010 | 1.0 |
| 540-59-0 | 1,2-Dichloroethene | 1.0 | U | 0.010 | 1.0 |
| 78-87-5 | 1,2-Dichloropropane | 1.0 | U | 0.010 | 1.0 |
| 541-73-1 | 1,3-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 106-46-7 | 1,4-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 78-93-3 | 2-Butanone | 5.0 | U | 0.010 | 5.0 |
| 501-78-6 | 2-Hexanone | 5.0 | U | 0.010 | 5.0 |
| 108-10-1 | 4-Methyl-2-pentanone | 5.0 | U | 0.010 | 5.0 |
| 67-64-1 | Acetone | 5.022 | 85 | 0.010 | 5.0 |
| 71-43-2 | Benzene | 1.0 | U | 0.010 | 1.0 |
| 75-27-4 | Bromodichloromethane | 1.0 | U | 0.010 | 1.0 |
| 75-25-2 | Bromoforn | 1.0 | U | 0.010 | 1.0 |
| 74-83-9 | Bromomethane | 1.0 | U | 0.010 | 1.0 |
| 75-15-0 | Carbon disulfide | 1.0 | U | 0.010 | 1.0 |
| 56-23-5 | Carbon tetrachloride | 1.0 | U | 0.010 | 1.0 |
| 106-90-7 | Chlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 75-00-3 | Chloroethane | 1.0 | U | 0.010 | 1.0 |
| 67-68-3 | Chloroform | 1.0 | U | 0.010 | 1.0 |
| 74-87-3 | Chloromethane | 1.0 | U | 0.010 | 1.0 |
| 124-48-1 | Dibromochloromethane | 1.0 | U | 0.010 | 1.0 |
| 10061-01-5 | cis-1,3-Dichloropropene | 1.0 | U | 0.010 | 1.0 |
| 10061-02-6 | trans-1,3-Dichloropropene | 1.0 | U | 0.010 | 1.0 |
| 100-41-4 | Ethylbenzene | 1.0 | U | 0.010 | 1.0 |
| 75-09-2 | Methylene chloride | 1.0218 | J | 0.010 | 2.0 |
| 100-42-5 | Styrene | 1.0 | U | 0.010 | 1.0 |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SKGW601009

Lab Name: GCAL Contract: _____

Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204031909

Matrix: (soil/water) Water

Sample wt/vol: 25 (g/ml) mL Lab Sample ID: 20403220804

Level: (low/med) _____ Lab File ID: 2040323/T2562

% Moisture: not dec. _____ Date Collected: 03/17/04 Time: 1115

GC Column: DB-624-30M ID: .53 (mm) Date Received: 03/19/04

Instrument ID: MSV2 Date Analyzed: 03/23/04 Time: 1505

Soil Extract Volume: _____ (µL) Dilution Factor: 1 Analyst: RSP

Soil Aliquot Volume: _____ (µL) Prep Batch: _____ Analytical Batch: 271108

CONCENTRATION UNITS: ug/L Analytical Method: OLCO 2.1

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|-----------|-------------------|--------|---|-------|-----|
| 127-18-4 | Tetrachloroethene | 1.0 | U | 0.010 | 1.0 |
| 108-88-3 | Toluene | 1.0 | U | 0.010 | 1.0 |
| 79-01-6 | Trichloroethene | 1.0 | U | 0.010 | 1.0 |
| 75-01-4 | Vinyl chloride | 1.0 | U | 0.010 | 1.0 |
| 1330-20-7 | Xylene (total) | 1.0 | U | 0.010 | 1.0 |

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

SKGW601009

Lab Name: GCAL Contract: _____
 Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204031909
 Matrix: Water Lab Sample ID: 20403220804
 Sample wt/vol: _____ Units: _____ Lab File ID: 2040323/T2562
 Level: (low/med) _____ Date Collected: 03/17/04 Time: 1115
 % Moisture: not dec. _____ Date Received: 03/19/04
 GC Column: D8-624-30M ID: .25 (mm) Date Analyzed: 03/23/04 Time: 1505
 Instrument ID: MSV2 Dilution Factor: 1 Analyst: RSP
 Soil Extract Volume: _____ (µL)
 Soil Aliquot Volume: _____ (µL)

Number TICs Found: 0

CONCENTRATION UNITS

| CAS NO. | COMPOUND | RT | EST. CONC. | Q |
|---------|------------------|----|------------|---|
| 1. | No tics detected | | | |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SKGW62A1009

Lab Name: GCAL

Contract: _____

Lab Code: LA024

Case No.: _____

SAS No.: _____

SDG No.: 204031909

Matrix: (soil/water) Water

Sample wt/vol: 25 (g/ml) mL

Lab Sample ID: 20403220805

Level: (low/med) _____

Lab File ID: 2040323/T2563

% Moisture: not dec. _____

Date Collected: 03/17/04

Time: 1150

GC Column: DB-624-30M ID: .53 (mm)

Date Received: 03/19/04

Instrument ID: MSV2

Date Analyzed: 03/23/04

Time: 1530

Soil Extract Volume: _____ (µL)

Dilution Factor: 1

Analyst: RSP

Soil Aliquot Volume: _____ (µL)

Prep Batch: _____

Analytical Batch: 271108

CONCENTRATION UNITS: ug/L

Analytical Method: OLCO 2.1

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|------------|---------------------------|--------|---|-------|-----|
| 71-55-6 | 1,1,1-Trichloroethane | 1.0 | U | 0.010 | 1.0 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1.0 | U | 0.010 | 1.0 |
| 79-00-5 | 1,1,2-Trichloroethane | 1.0 | U | 0.010 | 1.0 |
| 75-34-3 | 1,1-Dichloroethane | 1.0 | U | 0.010 | 1.0 |
| 75-35-4 | 1,1-Dichloroethene | 1.0 | U | 0.010 | 1.0 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 106-93-4 | 1,2-Dibromoethane | 1.0 | U | 0.010 | 1.0 |
| 95-50-1 | 1,2-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 107-06-2 | 1,2-Dichloroethane | 1.0 | U | 0.010 | 1.0 |
| 540-59-0 | 1,2-Dichloroethene | 1.0 | U | 0.010 | 1.0 |
| 78-87-5 | 1,2-Dichloropropane | 1.0 | U | 0.010 | 1.0 |
| 541-73-1 | 1,3-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 106-46-7 | 1,4-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 78-93-3 | 2-Butanone | 5.0 | U | 0.010 | 5.0 |
| 591-78-6 | 2-Hexanone | 5.0 | U | 0.010 | 5.0 |
| 108-10-1 | 4-Methyl-2-pentanone | 5.0 | U | 0.010 | 5.0 |
| 67-64-1 | Acetone | 5.0 | U | 0.010 | 5.0 |
| 71-43-2 | Benzene | 1.0 | U | 0.010 | 1.0 |
| 75-27-4 | Bromodichloromethane | 1.0 | U | 0.010 | 1.0 |
| 75-25-2 | Bromochloromethane | 1.0 | U | 0.010 | 1.0 |
| 74-83-9 | Bromomethane | 1.0 | U | 0.010 | 1.0 |
| 75-15-0 | Carbon disulfide | 1.0 | U | 0.010 | 1.0 |
| 56-23-5 | Carbon tetrachloride | 1.0 | U | 0.010 | 1.0 |
| 108-90-7 | Chlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 75-00-3 | Chloroethane | 1.0 | U | 0.010 | 1.0 |
| 67-66-3 | Chloroform | 1.0 | U | 0.010 | 1.0 |
| 74-87-3 | Chloromethane | 1.0 | U | 0.010 | 1.0 |
| 124-48-1 | Dibromochloromethane | 1.0 | U | 0.010 | 1.0 |
| 10061-01-5 | cis-1,3-Dichloropropene | 1.0 | U | 0.010 | 1.0 |
| 10061-02-6 | trans-1,3-Dichloropropene | 1.0 | U | 0.010 | 1.0 |
| 100-41-4 | Ethylbenzene | 1.0 | U | 0.010 | 1.0 |
| 75-09-2 | Methylene chloride | 2.0 | J | 0.010 | 2.0 |
| 100-42-5 | Styrene | 1.0 | U | 0.010 | 1.0 |

VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SKGW62A1009

Lab Name: GCAL

Contract:

Lab Code: LA024

Case No.:

SAS No.:

SDG No.: 204031909

Matrix: (soil/water) Water

Sample wt/vol: 25 (g/ml) mL

Lab Sample ID: 20403220805

Level: (low/med)

Lab File ID: 2040323/T2563

% Moisture: not dec.

Date Collected: 03/17/04

Time: 1150

GC Column: DB-624-30M

ID: 53

(mm)

Date Received: 03/19/04

Instrument ID: MSV2

Date Analyzed: 03/23/04

Time: 1530

Soil Extract Volume: (µL)

Dilution Factor: 1

Analyst: RSP

Soil Aliquot Volume: (µL)

Prep Batch:

Analytical Batch: 271108

CONCENTRATION UNITS ug/L

Analytical Method: OLCO 2.1

| GAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|-----------|-------------------|--------|---|-------|-----|
| 127-18-4 | Tetrachloroethene | < C | U | 0.010 | 1.0 |
| 106-88-3 | Toluene | < C | U | 0.010 | 1.0 |
| 79-01-6 | Trichloroethene | < C | U | 0.010 | 1.0 |
| 75-01-4 | Vinyl chloride | < C | U | 0.010 | 1.0 |
| 1330-20-7 | Xylene (total) | < C | U | 0.010 | 1.0 |

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

SKGW62A1009

Lab Name: GCAL Contract: _____

Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204031909

Matrix: Water Lab Sample ID: 20403220805

Sample wt/vol: _____ Units: _____ Lab File ID: 2040323/T2563

Level: (low/med) _____ Date Collected: 03/17/04 Time: 1150

% Moisture: not dec. _____ Date Received: 03/19/04

GC Column: DB-624-30M ID: .25 (mm) Date Analyzed: 03/23/04 Time: 1530

Instrument ID: MSV2 Dilution Factor: 1 Analyst: RSP

Soil Extract Volume: _____ (μ L)

Soil Aliquot Volume: _____ (μ L)

Number TICs Found: 0

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | RT | EST. CONC. | Q |
|---------|------------------|----|------------|---|
| 1. | No tics detected | | | |

VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SKGW641009

Lab Name: GCAL

Contract:

Lab Code: LA024

Case No.:

SAS No.:

SDG No.: 204031909

Matrix: (soil/water) Water

Sample wt/vol: 25 (g/m) mL

Lab Sample ID: 20403220806

Level: (low/med)

Lab File ID: 2040323/T2564

% Moisture: not dec.

Date Collected: 03/17/04

Time: 1410

GC Column: DB-624-30M ID: .53 (mm)

Date Received: 03/19/04

Instrument ID: MSV2

Date Analyzed: 03/23/04

Time: 1554

Soil Extract Volume: (µL)

Dilution Factor: 1

Analyst: RSP

Soil Aliquot Volume: (µL)

Prep Batch:

Analytical Batch: 271108

CONCENTRATION UNITS µg/L

Analytical Method: OLCO 2.1

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|------------|---------------------------|---------|---|-------|-----|
| 71-55-6 | 1,1,1-Trichloroethane | ND | U | 0.010 | 1.0 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | U | 0.010 | 1.0 |
| 79-00-5 | 1,1,2-Trichloroethane | ND | U | 0.010 | 1.0 |
| 75-34-3 | 1,1-Dichloroethane | ND | U | 0.010 | 1.0 |
| 75-35-4 | 1,1-Dichloroethene | ND | U | 0.010 | 1.0 |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | U | 0.010 | 1.0 |
| 106-93-4 | 1,2-Dibromoethane | ND | U | 0.010 | 1.0 |
| 95-50-1 | 1,2-Dichlorobenzene | ND | U | 0.010 | 1.0 |
| 107-06-2 | 1,2-Dichloroethane | ND | U | 0.010 | 1.0 |
| 540-59-0 | 1,2-Dichloroethene | ND | U | 0.010 | 1.0 |
| 78-87-5 | 1,2-Dichloropropane | ND | U | 0.010 | 1.0 |
| 541-73-1 | 1,3-Dichlorobenzene | ND | U | 0.010 | 1.0 |
| 106-46-7 | 1,4-Dichlorobenzene | ND | U | 0.010 | 1.0 |
| 78-93-3 | 2-Butanone | 5.0 | U | 0.010 | 5.0 |
| 591-78-6 | 2-Hexanone | 5.0 | U | 0.010 | 5.0 |
| 106-10-1 | 4-Methyl-2-pentanone | 5.0 | U | 0.010 | 5.0 |
| 67-64-1 | Acetone | 5.0 | U | 0.010 | 5.0 |
| 71-43-2 | Benzene | ND | U | 0.010 | 1.0 |
| 75-27-4 | Bromochloromethane | ND | U | 0.010 | 1.0 |
| 75-25-2 | Bromoforn | ND | U | 0.010 | 1.0 |
| 74-83-9 | Bromomethane | ND | U | 0.010 | 1.0 |
| 75-15-0 | Carbon disulfide | ND | U | 0.010 | 1.0 |
| 56-23-5 | Carbon tetrachloride | ND | U | 0.010 | 1.0 |
| 106-90-7 | Chlorobenzene | ND | U | 0.010 | 1.0 |
| 75-00-3 | Chloroethane | ND | U | 0.010 | 1.0 |
| 67-66-3 | Chloroform | ND | U | 0.010 | 1.0 |
| 74-87-3 | Chloromethane | ND | U | 0.010 | 1.0 |
| 124-48-1 | Dibromochloromethane | ND | U | 0.010 | 1.0 |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | U | 0.010 | 1.0 |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | U | 0.010 | 1.0 |
| 100-41-4 | Ethylbenzene | ND | U | 0.010 | 1.0 |
| 75-09-2 | Methylene chloride | 2.0.045 | J | 0.010 | 2.0 |
| 100-42-5 | Styrene | ND | U | 0.010 | 1.0 |

5/12/04
MPC

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SKGW641009

Lab Name: GCAL Contract: _____
 Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204031909
 Matrix: (soil/water) Water
 Sample wt/vol: 25 (g/ml) mL Lab Sample ID: 20403220806
 Level: (low/med) _____ Lab File ID: 2040323/T2564
 % Moisture: not dec. _____ Date Collected: 03/17/04 Time: 1410
 GC Column: DB-624-30M ID: .53 (mm) Date Received: 03/19/04
 Instrument ID: MSV2 Date Analyzed: 03/23/04 Time: 1554
 Soil Extract Volume: _____ (µL) Dilution Factor: 1 Analyst: RSP
 Soil Aliquot Volume: _____ (µL) Prep Batch: _____ Analytical Batch: 271108

CONCENTRATION UNITS: ug/L

Analytical Method: OLCO 2.1

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|-----------|-------------------|--------|---|-------|-----|
| 127-18-4 | Tetrachloroethene | 1.0 | U | 0.010 | 1.0 |
| 108-88-3 | Toluene | 1.0 | U | 0.010 | 1.0 |
| 79-01-6 | Trichloroethene | 1.0 | U | 0.010 | 1.0 |
| 75-01-4 | Vinyl chloride | 1.0 | U | 0.010 | 1.0 |
| 1330-20-7 | Xylene (total) | 1.0 | U | 0.010 | 1.0 |

VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

SKGW641009

Lab Name: GCAL Contract: _____
 Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204031909
 Matrix: Water Lab Sample ID: 20403220606
 Sample wt/vol: _____ Units: _____ Lab File ID: 2040323/T2564
 Level: (low/med) _____ Date Collected: 03/17/04 Time: 1410
 % Moisture: not dec. _____ Date Received: 03/19/04
 GC Column: DB-624-30M ID: 25 (mm) Date Analyzed: 03/23/04 Time: 1554
 Instrument ID: MSV2 Dilution Factor: 1 Analyst: RSP
 Soil Extract Volume: _____ (µL)
 Soil Aliquot Volume: _____ (µL)

Number TICs Found: 3

CONCENTRATION UNITS

| CAS NO. | COMPOUND | RT | EST. CONC. | Q |
|---------|------------------|----|------------|---|
| 1. | No pics detected | | | |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SKGW651009

Lab Name: GCAL Contract: _____

Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204031909

Matrix: (soil/water) Water

Sample wt/vol: 25 (g/ml) mL Lab Sample ID: 20403220807

Level: (low/med) _____ Lab File ID: 2040323/T2565

% Moisture: not dec. _____ Date Collected: 03/17/04 Time: 1440

GC Column: DB-624-30M ID: .53 (mm) Date Received: 03/19/04

Instrument ID: MSV2 Date Analyzed: 03/23/04 Time: 1619

Soil Extract Volume: _____ (μL) Dilution Factor: 1 Analyst: RSP

Soil Aliquot Volume: _____ (μL) Prep Batch: _____ Analytical Batch: 271108

CONCENTRATION UNITS: ug/L

Analytical Method: OLCO 2.1

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|------------|---------------------------|---------|---|-------|-----|
| 71-55-6 | 1,1,1-Trichloroethane | 1.0 | U | 0.010 | 1.0 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1.0 | U | 0.010 | 1.0 |
| 79-00-5 | 1,1,2-Trichloroethane | 1.0 | U | 0.010 | 1.0 |
| 75-34-3 | 1,1-Dichloroethane | 1.0 | U | 0.010 | 1.0 |
| 75-35-4 | 1,1-Dichloroethene | 1.0 | U | 0.010 | 1.0 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 106-83-4 | 1,2-Dibromoethane | 1.0 | U | 0.010 | 1.0 |
| 95-50-1 | 1,2-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 107-06-2 | 1,2-Dichloroethane | 1.0 | U | 0.010 | 1.0 |
| 540-59-0 | 1,2-Dichloroethene | 1.0 | U | 0.010 | 1.0 |
| 78-87-5 | 1,2-Dichloropropane | 1.0 | U | 0.010 | 1.0 |
| 541-73-1 | 1,3-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 106-46-7 | 1,4-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 78-93-3 | 2-Butanone | 5.0 | U | 0.010 | 5.0 |
| 591-78-6 | 2-Hexanone | 5.0 | U | 0.010 | 5.0 |
| 108-10-1 | 4-Methyl-2-pentanone | 5.0 | U | 0.010 | 5.0 |
| 67-64-1 | Acetone | 5.0 | U | 0.010 | 5.0 |
| 71-43-2 | Benzene | 1.0 | U | 0.010 | 1.0 |
| 75-27-4 | Bromodichloromethane | 1.0 | U | 0.010 | 1.0 |
| 75-25-2 | Bromoform | 1.0 | U | 0.010 | 1.0 |
| 74-83-9 | Bromomethane | 1.0 | U | 0.010 | 1.0 |
| 75-15-0 | Carbon disulfide | 1.0 | U | 0.010 | 1.0 |
| 56-23-5 | Carbon tetrachloride | 1.0 | U | 0.010 | 1.0 |
| 108-90-7 | Chlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 75-00-3 | Chloroethane | 1.0 | U | 0.010 | 1.0 |
| 67-66-3 | Chloroform | 1.0 | U | 0.010 | 1.0 |
| 74-87-3 | Chloromethane | 1.0 | U | 0.010 | 1.0 |
| 124-48-1 | Dibromochloromethane | 1.0 | U | 0.010 | 1.0 |
| 10061-01-5 | cis-1,3-Dichloropropene | 1.0 | U | 0.010 | 1.0 |
| 10061-02-6 | trans-1,3-Dichloropropene | 1.0 | U | 0.010 | 1.0 |
| 100-41-4 | Ethylbenzene | 1.0 | U | 0.010 | 1.0 |
| 75-09-2 | Methylene chloride | 2.0-2.1 | J | 0.010 | 2.0 |
| 100-42-5 | Styrene | 1.0 | U | 0.010 | 1.0 |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SKGW651009

Lab Name: GCAL

Contract:

Lab Code: LA024

Case No.:

SAS No

SDG No: 204031909

Matrix: (soil/water) Water

Sample wt/vol: 25 (g/m) mL

Lab Sample ID: 20403220607

Level: (low/mid)

Lab File ID: 2040323/T2565

% Moisture: not dec.

Date Collected: 03/17/04

Time: 1440

GC Column: DB-624-30M ID: .53 (mm)

Date Received: 03/19/04

Instrument ID: MSV2

Date Analyzed: 03/23/04

Time: 1619

Soil Extract Volume: (µL)

Dilution Factor: 1

Analyst: RSP

Soil Aliquot Volume: (µL)

Prep Batch

Analytical Batch: 271108

CONCENTRATION UNITS µg/L

Analytical Method: OLCO 2.1

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|-----------|-------------------|--------|---|-------|-----|
| 127-18-4 | Tetrachloroethene | 1.0 | U | 0.010 | 1.0 |
| 108-88-3 | Toluene | 1.0 | U | 0.010 | 1.0 |
| 79-01-6 | Trichloroethene | 1.0 | U | 0.010 | 1.0 |
| 75-01-4 | Vinyl chloride | 1.0 | U | 0.010 | 1.0 |
| 1330-20-7 | Xylene (total) | 1.0 | U | 0.010 | 1.0 |

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

SKGW651009

Lab Name: GCAL Contract: _____

Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204031909

Matrix: Water Lab Sample ID: 20403220807

Sample wt/vol: _____ Units: _____ Lab File ID: 2040323/T2565

Level: (low/med) _____ Date Collected: 03/17/04 Time: 1440

% Moisture: not dec. _____ Date Received: 03/19/04

GC Column: DB-624-30M ID: .25 (mm) Date Analyzed: 03/23/04 Time: 1619

Instrument ID: MSV2 Dilution Factor: 1 Analyst: RSP

Soil Extract Volume: _____ (µL)

Soil Aliquot Volume: _____ (µL)

Number TICs Found: 0

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | RT | EST. CONC. | Q |
|---------|------------------|----|------------|---|
| 1. | No tics detected | | | |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

TRIP BLANK

Lab Name: GCAL Contract: _____

Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204031909

Matrix: (soil/water) Water

Sample wt/vol: 25 (g/ml) 1 mL Lab Sample ID: 20403220814

Level: (low/med) _____ Lab File ID: 2040323/T2560

% Moisture: not dec. _____ Date Collected: 03/17/04 Time: 0000

GC Column: DB-624-30M ID: .53 (mm) Date Received: 03/19/04

Instrument ID: MSV2 Date Analyzed: 03/23/04 Time: 1416

Soil Extract Volume: _____ (µL) Dilution Factor: 1 Analyst: RSP

Soil Aliquot Volume: _____ (µL) Prep Batch: _____ Analytical Batch: 271108

Analytical Method: OLCO 2.1

CONCENTRATION UNITS: ug/L

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|------------|---------------------------|--------|---|-------|-----|
| 71-55-6 | 1,1,1-Trichloroethane | 1.0 | U | 0.010 | 1.0 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1.0 | U | 0.010 | 1.0 |
| 79-00-5 | 1,1,2-Trichloroethane | 1.0 | U | 0.010 | 1.0 |
| 75-34-3 | 1,1-Dichloroethane | 1.0 | U | 0.010 | 1.0 |
| 75-35-4 | 1,1-Dichloroethene | 1.0 | U | 0.010 | 1.0 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 106-83-4 | 1,2-Dibromoethane | 1.0 | U | 0.010 | 1.0 |
| 85-50-1 | 1,2-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 107-06-2 | 1,2-Dichloroethane | 1.0 | U | 0.010 | 1.0 |
| 540-58-0 | 1,2-Dichloroethene | 1.0 | U | 0.010 | 1.0 |
| 78-67-5 | 1,2-Dichloropropane | 1.0 | U | 0.010 | 1.0 |
| 541-73-1 | 1,3-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 106-46-7 | 1,4-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 78-83-3 | 2-Butanone | 5.0 | U | 0.010 | 5.0 |
| 591-78-6 | 2-Hexanone | 5.0 | U | 0.010 | 5.0 |
| 108-10-1 | 4-Methyl-2-pentanone | 5.0 | U | 0.010 | 5.0 |
| 67-64-1 | Acetone | 3.1 | J | 0.010 | 5.0 |
| 71-43-2 | Benzene | 1.0 | U | 0.010 | 1.0 |
| 75-27-4 | Bromodichloromethane | 1.0 | U | 0.010 | 1.0 |
| 75-25-2 | Bromoform | 1.0 | U | 0.010 | 1.0 |
| 74-83-9 | Bromomethane | 1.0 | U | 0.010 | 1.0 |
| 75-15-0 | Carbon disulfide | 1.0 | U | 0.010 | 1.0 |
| 56-23-5 | Carbon tetrachloride | 1.0 | U | 0.010 | 1.0 |
| 106-90-7 | Chlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 75-00-3 | Chloroethane | 1.0 | U | 0.010 | 1.0 |
| 67-66-3 | Chloroform | 1.0 | U | 0.010 | 1.0 |
| 74-87-3 | Chloromethane | 1.0 | U | 0.010 | 1.0 |
| 124-48-1 | Dibromochloromethane | 1.0 | U | 0.010 | 1.0 |
| 10061-01-5 | cis-1,3-Dichloropropene | 1.0 | U | 0.010 | 1.0 |
| 10061-02-6 | trans-1,3-Dichloropropene | 1.0 | U | 0.010 | 1.0 |
| 100-41-4 | Ethylbenzene | 1.0 | U | 0.010 | 1.0 |
| 75-09-2 | Methylene chloride | 2.0 | J | 0.010 | 2.0 |
| 100-42-5 | Styrene | 1.0 | U | 0.010 | 1.0 |

5/22/04
m

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

TRIP BLANK

Lab Name: GCAL Contract: _____

Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204031909

Matrix: (soil/water) Water

Sample wt/vol: 25 (g/ml) mL Lab Sample ID: 20403220814

Level: (low/med) _____ Lab File ID: 2040323/T2560

% Moisture: not dec. _____ Date Collected: 03/17/04 Time: 0000

GC Column: DB-624-30M ID: .53 (mm) Date Received: 03/19/04

Instrument ID: MSV2 Date Analyzed: 03/23/04 Time: 1416

Soil Extract Volume: _____ (μL) Dilution Factor: 1 Analyst: RSP

Soil Aliquot Volume: _____ (μL) Prep Batch: _____ Analytical Batch: 271108

Analytical Method: OLCO 2.1

CONCENTRATION UNITS: ug/L

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|-----------|-------------------|--------|---|-------|-----|
| 127-18-4 | Tetrachloroethene | 1.0 | U | 0.010 | 1.0 |
| 108-88-3 | Toluene | 1.0 | U | 0.010 | 1.0 |
| 79-01-6 | Trichloroethene | 1.0 | U | 0.010 | 1.0 |
| 75-01-4 | Vinyl chloride | 1.0 | U | 0.010 | 1.0 |
| 1330-20-7 | Xylene (total) | 1.0 | U | 0.010 | 1.0 |

VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

TRIP BLANK

Lab Name: GCAL Contract: _____

Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204031909

Matrix: Water Lab Sample ID: 20403220814

Sample w/vol: _____ Units: _____ Lab File ID: 2040323/T2560

Level: (low/med) _____ Date Collected: 03/17/04 Time: 0000

% Moisture: not dec. _____ Date Received: 03/19/04

GC Column: DB-624-30M ID: .53 (mm) Date Analyzed: 03/23/04 Time: 1416

Instrument ID: MSV2 Dilution Factor: 1 Analyst: RSP

Soil Extract Volume: _____ (µL)

Soil Aliquot Volume: _____ (µL)

Number TICs Found: 1

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | RT | EST. CONC. | Q |
|--------------|----------------|-------|------------|---|
| 1. 7446-09-5 | Sulfur dioxide | 1.412 | 1.92 | |

SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

Lab Name: GCAL Sample ID: SKGW581009
 Lab Code: LA024 Case No.: _____ Contract: _____
 SAS No.: _____ SDG No.: 204031909 Lab File ID: 2040326/S5023
 Matrix: Water Lab Sample ID: 20403190901
 Sample wt/vol: 1000 Units: mL Date Collected: 03/18/04 Time: 1148
 Level: (low/med) _____ Date Received: 03/19/04
 % Moisture: _____ decanted: (Y/N) _____ Date Extracted: 3/22/04
 GC Column: DB-5MS-30M ID: .25 (mm) Date Analyzed: 03/26/04 Time: 1612
 Concentrated Extract Volume: 1000 (µL) Dilution Factor: 1 Analyst: JAR3
 Injection Volume: 1.0 (µL) Prep Method: _____
 GPC Cleanup: (Y/N) N pH: _____ Analytical Method: OLMO 4.2
 Instrument ID: MSSV2

CONCENTRATION UNITS: ug/LPrep Batch: 271164 Analytical Batch: 271781

| CAS NO. | COMPOUND | RESULT | Q | MDL | PQL |
|-----------|-----------------------------|--------|---|-------|------|
| 95-95-4 | 2,4,5-Trichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 88-06-2 | 2,4,6-Trichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 120-83-2 | 2,4-Dichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 51-28-5 | 2,4-Dinitrophenol | 25.0 | U | 0.010 | 25.0 |
| 121-14-2 | 2,4-Dinitrotoluene | 10.0 | U | 0.010 | 10.0 |
| 606-20-2 | 2,6-Dinitrotoluene | 10.0 | U | 0.010 | 10.0 |
| 91-58-7 | 2-Chloronaphthalene | 10.0 | U | 0.010 | 10.0 |
| 95-57-8 | 2-Chlorophenol | 10.0 | U | 0.010 | 10.0 |
| 91-57-6 | 2-Methylnaphthalene | 10.0 | U | 0.010 | 10.0 |
| 83-74-4 | 2-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 83-75-5 | 2-Nitrophenol | 10.0 | U | 0.010 | 10.0 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 10.0 | U | 0.010 | 10.0 |
| 99-09-2 | 3-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 534-52-1 | 2-Methyl-4,6-dinitrophenol | 25.0 | U | 0.010 | 25.0 |
| 59-50-7 | 4-Chloro-3-methylphenol | 10.0 | U | 0.010 | 10.0 |
| 106-47-8 | 4-Chloroaniline | 10.0 | U | 0.010 | 10.0 |
| 7005-72-3 | 4-Chlorophenyl-phenylether | 10.0 | U | 0.010 | 10.0 |
| 106-44-5 | 4-Methylphenol (p-Cresol) | 10.0 | U | 0.010 | 10.0 |
| 83-32-9 | Acenaphthene | 10.0 | U | 0.010 | 10.0 |
| 208-96-8 | Acenaphthylene | 10.0 | U | 0.010 | 10.0 |
| 120-12-7 | Anthracene | 10.0 | U | 0.010 | 10.0 |
| 56-55-3 | Benzo(a)anthracene | 10.0 | U | 0.010 | 10.0 |
| 50-32-8 | Benzo(a)pyrene | 10.0 | U | 0.010 | 10.0 |
| 205-99-2 | Benzo(b)fluoranthene | 10.0 | U | 0.010 | 10.0 |
| 191-24-2 | Benzo(g,h,i)perylene | 10.0 | U | 0.010 | 10.0 |
| 207-08-9 | Benzo(k)fluoranthene | 10.0 | U | 0.010 | 10.0 |
| 111-91-1 | Bis(2-Chloroethoxy)methane | 10.0 | U | 0.010 | 10.0 |
| 111-44-4 | Bis(2-Chloroethyl)ether | 10.0 | U | 0.010 | 10.0 |
| 108-60-1 | bis(2-Chloroisopropyl)ether | 10.0 | U | 0.010 | 10.0 |
| 117-81-7 | bis(2-ethylhexyl)phthalate | 10.0 | U | 0.010 | 10.0 |

SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

Lab Name: GCAL Sample ID: SKGW581009
 Lab Code: LA024 Case No.: _____ Contract: _____
 SAS No.: _____ SDG No.: 204031909 Lab File ID: 2040326/S5023
 Matrix: Water Lab Sample ID: 20403190901
 Sample wt/vol: 1000 Units: mL Date Collected: 03/18/04 Time: 1148
 Level: (low/med) _____ Date Received: 03/19/04
 % Moisture: _____ decanted: (Y/N) _____ Date Extracted: 3/22/04
 GC Column: DB-5MS-30M ID: 25 (mm) Date Analyzed: 03/26/04 Time: 1612
 Concentrated Extract Volume: 1000 (µL) Dilution Factor: 1 Analyst: JAR3
 Injection Volume: 1.0 (µL) Prep Method: _____
 GPC Cleanup: (Y/N) N pH: _____ Analytical Method: OLMO 4.2
 Instrument ID: MSSV2
 Prep Batch: 271164 Analytical Batch: 271781

CONCENTRATION UNITS: µg/L

| CAS NO. | COMPOUND | RESULT | Q | MDL | PQL |
|----------|---------------------------|--------|---|-------|------|
| 101-55-3 | 4-Bromophenyl-phenylether | 10.0 | U | 0.010 | 10.0 |
| 85-88-7 | Butylbenzylthiathalate | 10.0 | U | 0.010 | 10.0 |
| 86-74-8 | Carbazole | 10.0 | U | 0.010 | 10.0 |
| 218-01-9 | Chrysene | 10.0 | U | 0.010 | 10.0 |
| 84-74-2 | Di-n-butylphthalate | 10.0 | U | 0.010 | 10.0 |
| 117-84-0 | Di-n-octylphthalate | 10.0 | U | 0.010 | 10.0 |
| 53-70-3 | Dibenz(a,h)anthracene | 10.0 | U | 0.010 | 10.0 |
| 132-64-9 | Dibenzofuran | 10.0 | U | 0.010 | 10.0 |
| 84-66-2 | Diethylphthalate | 10.0 | U | 0.010 | 10.0 |
| 131-11-3 | Dimethyl-phthalate | 10.0 | U | 0.010 | 10.0 |
| 105-67-9 | 2,4-Dimethylphenol | 10.0 | U | 0.010 | 10.0 |
| 206-44-0 | Fluoranthene | 10.0 | U | 0.010 | 10.0 |
| 86-73-7 | Fluorene | 10.0 | U | 0.010 | 10.0 |
| 118-74-1 | Hexachlorobenzene | 10.0 | U | 0.010 | 10.0 |
| 87-68-3 | Hexachlorobutadiene | 10.0 | U | 0.010 | 10.0 |
| 77-47-4 | Hexachlorocyclopentadiene | 10.0 | U | 0.010 | 10.0 |
| 67-72-1 | Hexachloroethane | 10.0 | U | 0.010 | 10.0 |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | 10.0 | U | 0.010 | 10.0 |
| 78-59-1 | Isophorone | 10.0 | U | 0.010 | 10.0 |
| 91-20-3 | Naphthalene | 10.0 | U | 0.010 | 10.0 |
| 100-01-6 | 4-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 98-95-3 | Nitrobenzene | 10.0 | U | 0.010 | 10.0 |
| 100-02-7 | 4-Nitrophenol | 25.0 | U | 0.010 | 25.0 |
| 87-86-5 | Pentachlorophenol | 25.0 | U | 0.010 | 25.0 |
| 85-01-8 | Phenanthrene | 10.0 | U | 0.010 | 10.0 |
| 108-95-2 | Phenol | 10.0 | U | 0.010 | 10.0 |
| 129-00-0 | Pyrene | 10.0 | U | 0.010 | 10.0 |
| 621-64-7 | N-Nitroso-α-n-propylamine | 10.0 | U | 0.010 | 10.0 |
| 86-30-6 | N-Nitrosodiphenylamine | 10.0 | U | 0.010 | 10.0 |
| 95-48-7 | o-Cresol | 10.0 | U | 0.010 | 10.0 |

1F
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

Lab Name: GCAL Sample ID: SKGW581009
Lab Code: LA024 2 Case No.: _____ Contract: _____
SAS No.: _____ SDG No.: 204031909 Lab File ID: S5023
Matrix: Water Lab Sample ID: 20403190901
Sample wt/vol: _____ Units: _____ Date Collected: 03/18/04 Time: 1148
Level: (low/med) _____ Date Received: 03/19/04
% Moisture: not dec. _____ Date Extracted: 3/22/04
GC Column: RTX-SMS-30 ID: .53 (mm) Date Analyzed: 03/26/04 Time: 1612
Concentrated Extract Volume: 1000 (µL) Dilution Factor: 1 Analyst: RLW
Injection Volume: 1.0 (µL) Prep Method: _____
GPC Cleanup: (Y/N) N pH: _____ Analytical Method: SW-846 8270C
Instrument ID: MSSV2

Number TICs Found : 0

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | RT | EST. CONC. | Q |
|---------|------------------|----|------------|---|
| 1. | No tics detected | | | |

5/17/04
msa

SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

Lab Name: GCAL
 Lab Code: LA024 Case No.: _____
 SAS No.: _____ SDG No.: 204031909
 Matrix: Water
 Sample wt/vol: 1000 Units mL
 Level: (low/med) _____
 % Moisture: _____ Decanted: (Y/N) _____
 GC Column: DB-5MS-30M ID: 25 (mm)
 Concentrated Extract Volume: 1000 (µL)
 Injection Volume: _____ (µL)
 GPC Cleanup: (Y/N) N pH: _____

Sample ID: SKGW58DUP1009
 Contract: _____
 Lab File ID: 2040326/S5024
 Lab Sample ID: 20403190902
 Date Collected: 03/18/04 Time: 1208
 Date Received: 03/19/04
 Date Extracted: 3/22/04
 Date Analyzed: 03/26/04 Time: 1640
 Dilution Factor: 1 Analyst: JAR3
 Prep Method: _____
 Analytical Method: OLMO 4.2
 Instrument ID: MSSV2

CONCENTRATION UNITS µg/L

Prep Batch: 271164 Analytical Batch: 271781

| CAS NO. | COMPOUND | RESULT | Q | MDL | PQL |
|-----------|-----------------------------|--------|---|-------|------|
| 95-95-4 | 2,4,5-Trichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 88-06-2 | 2,4,6-Trichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 120-83-2 | 2,4-Dichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 51-28-5 | 2,4-Dinitrochloro | 25.0 | U | 0.010 | 25.0 |
| 121-14-2 | 2,4-Dinitrotoluene | 10.0 | U | 0.010 | 10.0 |
| 606-20-2 | 2,6-Dinitrotoluene | 10.0 | U | 0.010 | 10.0 |
| 91-58-7 | 2-Chloronaphthalene | 10.0 | U | 0.010 | 10.0 |
| 95-57-8 | 2-Chlorophenol | 10.0 | U | 0.010 | 10.0 |
| 91-57-6 | 2-Methylnaphthalene | 10.0 | U | 0.010 | 10.0 |
| 88-74-4 | 2-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 88-75-5 | 2-Nitrophenol | 10.0 | U | 0.010 | 10.0 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 10.0 | U | 0.010 | 10.0 |
| 99-09-2 | 3-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 534-52-1 | 2-Methyl-4,6-dinitrophenol | 25.0 | U | 0.010 | 25.0 |
| 59-50-7 | 4-Chloro-3-methylphenol | 10.0 | U | 0.010 | 10.0 |
| 108-47-8 | 4-Chloroaniline | 10.0 | U | 0.010 | 10.0 |
| 7005-72-3 | 4-Chlorophenyl-phenylether | 10.0 | U | 0.010 | 10.0 |
| 108-44-5 | 4-Methylphenol (p-Cresol) | 10.0 | U | 0.010 | 10.0 |
| 83-32-9 | Acenaphthene | 10.0 | U | 0.010 | 10.0 |
| 208-96-8 | Acenaphthylene | 10.0 | U | 0.010 | 10.0 |
| 120-12-7 | Anthracene | 10.0 | U | 0.010 | 10.0 |
| 56-55-3 | Benzoxanthracene | 10.0 | U | 0.010 | 10.0 |
| 50-32-8 | Benzofluoranthene | 10.0 | U | 0.010 | 10.0 |
| 205-99-2 | Benzofluoranthene | 10.0 | U | 0.010 | 10.0 |
| 191-24-2 | Benzofluoranthene | 10.0 | U | 0.010 | 10.0 |
| 207-08-9 | Benzofluoranthene | 10.0 | U | 0.010 | 10.0 |
| 111-91-1 | Bis(2-Chloroethoxy)methane | 10.0 | U | 0.010 | 10.0 |
| 111-44-4 | Bis(2-Chloroethyl)ether | 10.0 | U | 0.010 | 10.0 |
| 108-60-1 | bis(2-Chloroisopropyl)ether | 10.0 | U | 0.010 | 10.0 |
| 117-81-7 | bis(2-ethoxyethyl)phthalate | 10.0 | U | 0.010 | 10.0 |

5/12/04
mm

SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

Lab Name: GCALSample ID: SKGW58DUP1009Lab Code: LA024 Case No.: _____

Contract: _____

SAS No.: _____ SDG No.: 204031909Lab File ID: 2040326/S5024Matrix: WaterLab Sample ID: 20403190902Sample wt/vol: 1000 Units: mLDate Collected: 03/18/04 Time: 1208

Level: (low/med) _____

Date Received: 03/19/04

% Moisture: _____ decanted: (Y/N) _____

Date Extracted: 3/22/04GC Column: DB-5MS-30M ID: .25 (mm)Date Analyzed: 03/26/04 Time: 1640Concentrated Extract Volume: 1000 (µL)Dilution Factor: 1 Analyst: JAR3Injection Volume: 1.0 (µL)

Prep Method: _____

GPC Cleanup: (Y/N) N pH: _____Analytical Method: OLMO 4.2Instrument ID: MSSV2CONCENTRATION UNITS: ug/LPrep Batch: 271164 Analytical Batch: 271781

| CAS NO. | COMPOUND | RESULT | Q | MDL | PQL |
|----------|----------------------------|--------|---|-------|------|
| 101-55-3 | 4-Bromophenyl-phenylether | 10.0 | U | 0.010 | 10.0 |
| 85-68-7 | Butylbenzylphthalate | 10.0 | U | 0.010 | 10.0 |
| 86-74-8 | Carbazole | 10.0 | U | 0.010 | 10.0 |
| 218-01-9 | Chrysene | 10.0 | U | 0.010 | 10.0 |
| 84-74-2 | Di-n-butylphthalate | 10.0 | U | 0.010 | 10.0 |
| 117-84-0 | Di-n-octylphthalate | 10.0 | U | 0.010 | 10.0 |
| 53-70-3 | Dibenz(a,h)anthracene | 10.0 | U | 0.010 | 10.0 |
| 132-64-9 | Dibenzofuran | 10.0 | U | 0.010 | 10.0 |
| 84-66-2 | Diethylphthalate | 10.0 | U | 0.010 | 10.0 |
| 131-11-3 | Dimethyl-phthalate | 10.0 | U | 0.010 | 10.0 |
| 105-67-9 | 2,4-Dimethylphenol | 10.0 | U | 0.010 | 10.0 |
| 206-44-0 | Fluoranthene | 10.0 | U | 0.010 | 10.0 |
| 86-73-7 | Fluorene | 10.0 | U | 0.010 | 10.0 |
| 118-74-1 | Hexachlorobenzene | 10.0 | U | 0.010 | 10.0 |
| 87-68-3 | Hexachlorobutadiene | 10.0 | U | 0.010 | 10.0 |
| 77-47-4 | Hexachlorocyclopentadiene | 10.0 | U | 0.010 | 10.0 |
| 67-72-1 | Hexachloroethane | 10.0 | U | 0.010 | 10.0 |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | 10.0 | U | 0.010 | 10.0 |
| 78-59-1 | Isophorone | 10.0 | U | 0.010 | 10.0 |
| 91-20-3 | Naphthalene | 10.0 | U | 0.010 | 10.0 |
| 100-01-6 | 4-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 98-95-3 | Nitrobenzene | 10.0 | U | 0.010 | 10.0 |
| 100-02-7 | 4-Nitrophenol | 25.0 | U | 0.010 | 25.0 |
| 87-86-5 | Pentachlorophenol | 25.0 | U | 0.010 | 25.0 |
| 85-01-8 | Phenanthrene | 10.0 | U | 0.010 | 10.0 |
| 108-95-2 | Phenol | 10.0 | U | 0.010 | 10.0 |
| 129-00-0 | Pyrene | 10.0 | U | 0.010 | 10.0 |
| 621-64-7 | N-Nitroso-di-n-propylamine | 10.0 | U | 0.010 | 10.0 |
| 86-30-6 | N-Nitrosodiphenylamine | 10.0 | U | 0.010 | 10.0 |
| 95-48-7 | o-Cresol | 10.0 | U | 0.010 | 10.0 |

4J
4J5/13/04
MK

SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

| | |
|---|---|
| Lab Name: <u>GCAL</u> | Sample ID: <u>SKGW58DUP1009</u> |
| Lab Code: <u>LA024</u> <u>2</u> Case No.: _____ | Contract: _____ |
| SAS No.: _____ SDG No.: <u>204031909</u> | Lab File ID: <u>S5024</u> |
| Matrix: <u>Water</u> | Lab Sample ID: <u>20403190902</u> |
| Sample wt/vol: _____ Units: _____ | Date Collected: <u>03/18/04</u> Time: <u>1208</u> |
| Level: (low/med) _____ | Date Received: <u>03/19/04</u> |
| % Moisture: not dec. _____ | Date Extracted: <u>3/22/04</u> |
| GC Column: <u>RTX-5MS-30</u> ID: <u>53</u> (mm) | Date Analyzed: <u>03/26/04</u> Time: <u>1640</u> |
| Concentrated Extract Volume: <u>1000</u> (µL) | Dilution Factor: <u>1</u> Analyst: <u>RLW</u> |
| Injection Volume: <u>1.0</u> (µL) | Prep Method: _____ |
| GPC Cleanup: (Y/N) <u>N</u> pH: _____ | Analytical Method: <u>SW-846 8270C</u> |
| | Instrument ID: <u>MSSV2</u> |

Number TICs Found: 0

CONCENTRATION UNITS

| CAS NO. | COMPOUND | RT | EST. CONC. | Q |
|---------|------------------|----|------------|---|
| 1. | No tics detected | | | |

5/17/07
MS

SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

Lab Name: GCAL Sample ID: SKGW631009
 Lab Code: LA024 Case No.: _____ Contract: _____
 SAS No.: _____ SDG No.: 204031909 Lab File ID: 2040326/S5025
 Matrix: Water Lab Sample ID: 20403190903
 Sample wt/vol: 1000 Units: mL Date Collected: 03/18/04 Time: 1430
 Level: (low/med) _____ Date Received: 03/19/04
 % Moisture: _____ decanted: (Y/N) _____ Date Extracted: 3/22/04
 GC Column: DB-5MS-30M ID: .25 (mm) Date Analyzed: 03/26/04 Time: 1707
 Concentrated Extract Volume: 1000 (µL) Dilution Factor: 1 Analyst: JAR3
 Injection Volume: 1.0 (µL) Prep Method: _____
 GPC Cleanup: (Y/N) N pH: _____ Analytical Method: OLMO 4.2
 Instrument ID: MSSV2

CONCENTRATION UNITS: ug/LPrep Batch: 271164 Analytical Batch: 271781

| CAS NO. | COMPOUND | RESULT | Q | MDL | PQL |
|-----------|-----------------------------|--------|---|-------|------|
| 95-95-4 | 2,4,5-Trichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 88-06-2 | 2,4,6-Trichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 120-83-2 | 2,4-Dichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 51-28-5 | 2,4-Dinitrophenol | 25.0 | U | 0.010 | 25.0 |
| 121-14-2 | 2,4-Dinitrotoluene | 10.0 | U | 0.010 | 10.0 |
| 606-20-2 | 2,6-Dinitrotoluene | 10.0 | U | 0.010 | 10.0 |
| 91-58-7 | 2-Chloronaphthalene | 10.0 | U | 0.010 | 10.0 |
| 95-57-8 | 2-Chlorophenol | 10.0 | U | 0.010 | 10.0 |
| 91-57-6 | 2-Methylnaphthalene | 10.0 | U | 0.010 | 10.0 |
| 88-74-4 | 2-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 88-75-5 | 2-Nitrophenol | 10.0 | U | 0.010 | 10.0 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 10.0 | U | 0.010 | 10.0 |
| 99-09-2 | 3-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 534-52-1 | 2-Methyl-4,6-dinitrophenol | 25.0 | U | 0.010 | 25.0 |
| 59-50-7 | 4-Chloro-3-methylphenol | 10.0 | U | 0.010 | 10.0 |
| 106-47-8 | 4-Chloroaniline | 10.0 | U | 0.010 | 10.0 |
| 7005-72-3 | 4-Chlorophenyl-phenylether | 10.0 | U | 0.010 | 10.0 |
| 106-44-5 | 4-Methylphenol (p-Cresol) | 10.0 | U | 0.010 | 10.0 |
| 83-32-9 | Acenaphthene | 10.0 | U | 0.010 | 10.0 |
| 208-96-8 | Acenaphthylene | 10.0 | U | 0.010 | 10.0 |
| 120-12-7 | Anthracene | 10.0 | U | 0.010 | 10.0 |
| 56-55-3 | Benzo(a)anthracene | 10.0 | U | 0.010 | 10.0 |
| 50-32-8 | Benzo(a)pyrene | 10.0 | U | 0.010 | 10.0 |
| 205-99-2 | Benzo(b)fluoranthene | 10.0 | U | 0.010 | 10.0 |
| 191-24-2 | Benzo(g,h,i)perylene | 10.0 | U | 0.010 | 10.0 |
| 207-08-9 | Benzo(k)fluoranthene | 10.0 | U | 0.010 | 10.0 |
| 111-91-1 | Bis(2-Chloroethoxy)methane | 10.0 | U | 0.010 | 10.0 |
| 111-44-4 | Bis(2-Chloroethyl)ether | 10.0 | U | 0.010 | 10.0 |
| 108-60-1 | bis(2-Chloroisopropyl)ether | 10.0 | U | 0.010 | 10.0 |
| 117-81-7 | bis(2-ethylhexyl)phthalate | 10.0 | U | 0.010 | 10.0 |

SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

Lab Name: GCAL
 Lab Code: LA024 Case No.: _____
 SAS No.: _____ SDG No.: 204031909
 Matrix: Water
 Sample wt/vol: 1000 Units: mL
 Level: (low/med) _____
 % Moisture: _____ decanted: (Y/N) _____
 GC Column: DB-5MS-30M ID: 25 (mm)
 Concentrated Extract Volume: 1000 (µL)
 Injection Volume: 1.0 (µL)
 GPC Cleanup: (Y/N) N pH: _____

Sample ID: SKGW631009
 Contract: _____
 Lab File ID: 2040326/S5025
 Lab Sample ID: 20403190903
 Date Collected: 03/18/04 Time: 1430
 Date Received: 03/19/04
 Date Extracted: 3/22/04
 Date Analyzed: 03/26/04 Time: 1707
 Dilution Factor: 1 Analyst: JAR3
 Prep Method: _____
 Analytical Method: OLMO 4.2
 Instrument ID: MSSV2

CONCENTRATION UNITS: µg/L

Prep Batch: 271164 Analytical Batch: 271781

| CAS NO. | COMPOUND | RESULT | Q | MDL | PQL |
|----------|---------------------------|--------|---|-------|------|
| 101-55-3 | 4-Bromophenyl-phenylether | 10.0 | U | 0.010 | 10.0 |
| 85-68-7 | Butylbenzophthalate | 0.610 | J | 0.010 | 10.0 |
| 86-74-8 | Carbazole | 10.0 | U | 0.010 | 10.0 |
| 218-01-9 | Chrysene | 10.0 | U | 0.010 | 10.0 |
| 84-74-2 | Di-n-butylphthalate | 10.0 | U | 0.010 | 10.0 |
| 117-84-0 | Di-n-octylphthalate | 10.0 | U | 0.010 | 10.0 |
| 53-70-3 | Dibenz(a,h)anthracene | 10.0 | U | 0.010 | 10.0 |
| 132-64-9 | Dibenzofuran | 10.0 | U | 0.010 | 10.0 |
| 84-66-2 | Diethylphthalate | 10.0 | U | 0.010 | 10.0 |
| 131-11-3 | Dimethyl-phthalate | 10.0 | U | 0.010 | 10.0 |
| 105-67-8 | 2,4-Dimethylphenol | 10.0 | U | 0.010 | 10.0 |
| 206-44-0 | Fluoranthene | 10.0 | U | 0.010 | 10.0 |
| 86-73-7 | Fluorene | 10.0 | U | 0.010 | 10.0 |
| 118-74-1 | Hexachlorobenzene | 10.0 | U | 0.010 | 10.0 |
| 87-68-3 | Hexachlorobutadiene | 10.0 | U | 0.010 | 10.0 |
| 77-47-4 | Hexachlorocyclopentadiene | 10.0 | U | 0.010 | 10.0 |
| 67-72-1 | Hexachloroethane | 10.0 | U | 0.010 | 10.0 |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | 10.0 | U | 0.010 | 10.0 |
| 78-59-1 | Isophorone | 10.0 | U | 0.010 | 10.0 |
| 91-20-3 | Naphthalene | 10.0 | U | 0.010 | 10.0 |
| 100-01-6 | 4-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 98-95-3 | Nitrobenzene | 10.0 | U | 0.010 | 10.0 |
| 100-02-7 | 4-Nitrophenol | 25.0 | U | 0.010 | 25.0 |
| 87-86-5 | Pentachlorononol | 25.0 | U | 0.010 | 25.0 |
| 85-01-8 | Phenanthrene | 10.0 | U | 0.010 | 10.0 |
| 108-95-2 | Phenol | 10.0 | U | 0.010 | 10.0 |
| 129-00-0 | Pyrene | 10.0 | U | 0.010 | 10.0 |
| 621-64-7 | N-Nitroso-d-n-propylamine | 10.0 | U | 0.010 | 10.0 |
| 86-30-6 | N-Nitrosodiphenylamine | 10.0 | U | 0.010 | 10.0 |
| 95-48-7 | o-Cresol | 10.0 | U | 0.010 | 10.0 |

SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

| | |
|--|---|
| Lab Name: <u>GCAL</u> | Sample ID: <u>SKGW631009</u> |
| Lab Code: <u>LA024</u> <u>2</u> Case No.: _____ | Contract: _____ |
| SAS No.: _____ SDG No.: <u>204031909</u> | Lab File ID: <u>S5025</u> |
| Matrix: <u>Water</u> | Lab Sample ID: <u>20403190903</u> |
| Sample wt/vol: _____ Units: _____ | Date Collected: <u>03/18/04</u> Time: <u>1430</u> |
| Level: (low/med) _____ | Date Received: <u>03/19/04</u> |
| % Moisture: not dec. _____ | Date Extracted: <u>3/22/04</u> |
| GC Column: <u>RTX-5MS-30</u> ID: <u>.53</u> (mm) | Date Analyzed: <u>03/26/04</u> Time: <u>1707</u> |
| Concentrated Extract Volume: <u>1000</u> (µL) | Dilution Factor: <u>1</u> Analyst: <u>RLW</u> |
| Injection Volume: <u>1.0</u> (µL) | Prep Method: _____ |
| GPC Cleanup: (Y/N) <u>N</u> pH: _____ | Analytical Method: <u>SW-846 8270C</u> |
| | Instrument ID: <u>MSSV2</u> |

Number TICs Found : 0

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | RT | EST. CONC. | Q |
|---------|------------------|----|------------|---|
| 1. | No tics detected | | | |

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SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

| | |
|---|---|
| Lab Name: <u>GCAL</u> | Sample ID: <u>SKGW611009</u> |
| Lab Code: <u>LA024</u> Case No.: _____ | Contract: _____ |
| SAS No.: _____ SDG No.: <u>204031909</u> | Lab File ID: <u>2040326/S5026</u> |
| Matrix: <u>Water</u> | Lab Sample ID: <u>20403190904</u> |
| Sample wt/vol: <u>1000</u> Units: <u>mL</u> | Date Collected: <u>03/18/04</u> Time: <u>1535</u> |
| Level: (low/med) _____ | Date Received: <u>03/19/04</u> |
| % Moisture: _____ decanted: (Y/N) _____ | Date Extracted: <u>3/22/04</u> |
| GC Column: <u>DB-SMS-30M</u> ID: <u>25</u> (mm) | Date Analyzed: <u>03/26/04</u> Time: <u>1735</u> |
| Concentrated Extract Volume: <u>1000</u> (µL) | Dilution Factor: <u>1</u> Analyst: <u>JAR3</u> |
| Injection Volume: <u>1.0</u> (µL) | Prep Method: _____ |
| GPC Cleanup: (Y/N) <u>N</u> pH: _____ | Analytical Method: <u>OLMO 4.2</u> |
| Instrument ID: <u>MSSV2</u> | |

CONCENTRATION UNITS: µg/L

CAS NO. COMPOUND

RESULT Q MDL PQL

| | | | | | |
|-----------|------------------------------|------|---|-------|------|
| 95-95-4 | 2,4,5-Trichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 88-06-2 | 2,4,6-Trichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 120-83-2 | 2,4-Dichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 51-28-5 | 2,4-Dinitrophenol | 25.0 | U | 0.010 | 25.0 |
| 121-14-2 | 2,4-Dinitrotoluene | 10.0 | U | 0.010 | 10.0 |
| 806-20-2 | 2,6-Dinitrotoluene | 10.0 | U | 0.010 | 10.0 |
| 91-58-7 | 2-Chloronaphthalene | 10.0 | U | 0.010 | 10.0 |
| 95-57-8 | 2-Chlorophenol | 10.0 | U | 0.010 | 10.0 |
| 91-57-6 | 2-Methylnaphthalene | 10.0 | U | 0.010 | 10.0 |
| 88-74-4 | 2-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 88-75-5 | 2-Nitrophenol | 10.0 | U | 0.010 | 10.0 |
| 91-84-1 | 3,3'-Dichlorodiphenylmethane | 10.0 | U | 0.010 | 10.0 |
| 99-09-2 | 3-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 534-52-1 | 2-Methyl-4,6-dinitrophenol | 25.0 | U | 0.010 | 25.0 |
| 59-50-7 | 4-Chloro-3-methylphenol | 10.0 | U | 0.010 | 10.0 |
| 106-47-8 | 4-Chloroaniline | 10.0 | U | 0.010 | 10.0 |
| 7005-72-3 | 4-Chlorophenyl-phenylether | 10.0 | U | 0.010 | 10.0 |
| 106-44-5 | 4-Methylphenol (p-Cresol) | 10.0 | U | 0.010 | 10.0 |
| 83-32-9 | Acenaphthene | 10.0 | U | 0.010 | 10.0 |
| 208-96-8 | Acenaphthylene | 10.0 | U | 0.010 | 10.0 |
| 120-12-7 | Anthracene | 10.0 | U | 0.010 | 10.0 |
| 56-55-3 | Benzo(a)anthracene | 10.0 | U | 0.010 | 10.0 |
| 50-32-8 | Benzo(a)pyrene | 10.0 | U | 0.010 | 10.0 |
| 205-99-2 | Benzo(b)fluoranthene | 10.0 | U | 0.010 | 10.0 |
| 191-24-2 | Benzo(g,h,i)perylene | 10.0 | U | 0.010 | 10.0 |
| 207-08-9 | Benzo(k)fluoranthene | 10.0 | U | 0.010 | 10.0 |
| 111-91-1 | Bis(2-Chloroethoxy)methane | 10.0 | U | 0.010 | 10.0 |
| 111-44-4 | Bis(2-Chloroethyl) ether | 10.0 | U | 0.010 | 10.0 |
| 106-60-1 | bis(2-Chloroisopropyl) ether | 10.0 | U | 0.010 | 10.0 |
| 117-81-7 | bis(2-ethoxy) phthalate | 10.0 | U | 0.010 | 10.0 |

SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

Lab Name: GCAL Sample ID: SKGW611009

Lab Code: LA024 Case No.: _____ Contract: _____

SAS No.: _____ SDG No.: 204031909 Lab File ID: 2040326/S5026

Matrix: Water Lab Sample ID: 20403190904

Sample wt/vol: 1000 Units: mL Date Collected: 03/18/04 Time: 1535

Level: (low/med) _____ Date Received: 03/19/04

% Moisture: _____ decanted: (Y/N) _____ Date Extracted: 3/22/04

GC Column: DB-5MS-30M ID: .25 (mm) Date Analyzed: 03/26/04 Time: 1735

Concentrated Extract Volume: 1000 (µL) Dilution Factor: 1 Analyst: JAR3

Injection Volume: 1.0 (µL) Prep Method: _____

GPC Cleanup: (Y/N) N pH: _____ Analytical Method: OLMO 4.2

Instrument ID: MSSV2

Prep Batch: 271164 Analytical Batch: 271781

CONCENTRATION UNITS: ug/L

| CAS NO. | COMPOUND | RESULT | Q | MDL | PQL |
|----------|----------------------------|--------|---|-------|------|
| 101-55-3 | 4-Bromophenyl-phenylether | 10.0 | U | 0.010 | 10.0 |
| 95-68-7 | Butylbenzylphthalate | 10.0 | U | 0.010 | 10.0 |
| 96-74-8 | Carbazole | 10.0 | U | 0.010 | 10.0 |
| 218-01-9 | Chrysene | 10.0 | U | 0.010 | 10.0 |
| 34-74-2 | Di-n-butylphthalate | 10.0 | U | 0.010 | 10.0 |
| 117-84-0 | Di-n-octylphthalate | 10.0 | U | 0.010 | 10.0 |
| 53-70-3 | Dibenz(a,h)anthracene | 10.0 | U | 0.010 | 10.0 |
| 132-64-9 | Dibenzofuran | 10.0 | U | 0.010 | 10.0 |
| 84-86-2 | Diethylphthalate | 10.0 | U | 0.010 | 10.0 |
| 131-11-3 | Dimethyl-phthalate | 10.0 | U | 0.010 | 10.0 |
| 105-67-9 | 2,4-Dimethylphenol | 10.0 | U | 0.010 | 10.0 |
| 206-44-0 | Fluoranthene | 10.0 | U | 0.010 | 10.0 |
| 86-73-7 | Fluorene | 10.0 | U | 0.010 | 10.0 |
| 118-74-1 | Hexachlorobenzene | 10.0 | U | 0.010 | 10.0 |
| 87-68-3 | Hexachlorobutadiene | 10.0 | U | 0.010 | 10.0 |
| 77-47-4 | Hexachlorocyclopentadiene | 10.0 | U | 0.010 | 10.0 |
| 67-72-1 | Hexachloroethane | 10.0 | U | 0.010 | 10.0 |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | 10.0 | U | 0.010 | 10.0 |
| 78-59-1 | Isophorone | 10.0 | U | 0.010 | 10.0 |
| 91-20-3 | Naphthalene | 10.0 | U | 0.010 | 10.0 |
| 100-01-6 | 4-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 98-95-3 | Nitrobenzene | 10.0 | U | 0.010 | 10.0 |
| 100-02-7 | 4-Nitrophenol | 25.0 | U | 0.010 | 25.0 |
| 87-86-5 | Pentachlorophenol | 25.0 | U | 0.010 | 25.0 |
| 85-01-8 | Phenanthrene | 10.0 | U | 0.010 | 10.0 |
| 108-95-2 | Phenol | 10.0 | U | 0.010 | 10.0 |
| 129-00-0 | Pyrene | 10.0 | U | 0.010 | 10.0 |
| 621-64-7 | N-Nitroso-di-n-propylamine | 10.0 | U | 0.010 | 10.0 |
| 86-30-6 | N-Nitrosodiphenylamine | 10.0 | U | 0.010 | 10.0 |
| 95-48-7 | o-Cresol | 10.0 | U | 0.010 | 10.0 |

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SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

| | |
|---|---|
| Lab Name: <u>GCAL</u> | Sample ID: <u>SKGW611009</u> |
| Lab Code: <u>LA024</u> <u>2</u> Case No.: _____ | Contract: _____ |
| SAS No.: _____ SDG No.: <u>204031909</u> | Lab File ID: <u>S5026</u> |
| Matrix: <u>Water</u> | Lab Sample ID: <u>20403190904</u> |
| Sample wt/vol: _____ Units _____ | Date Collected: <u>03/18/04</u> Time: <u>1535</u> |
| Level: (low/med) _____ | Date Received: <u>03/19/04</u> |
| % Moisture: not dec. _____ | Date Extracted: <u>3/22/04</u> |
| GC Column: <u>RTX-5MS-30</u> ID: <u>53</u> (mm) | Date Analyzed: <u>03/26/04</u> Time: <u>1735</u> |
| Concentrated Extract Volume: <u>1000</u> (µL) | Dilution Factor: <u>1</u> Analyst: <u>RLW</u> |
| Injection Volume: <u>10</u> (µL) | Prep Method: _____ |
| GPC Cleanup: (Y/N) <u>N</u> pH: _____ | Analytical Method: <u>SW-846 8270C</u> |
| | Instrument ID: <u>MSSV2</u> |

Number TICs Found : 0

CONCENTRATION UNITS

| CAS NO. | COMPOUND | RT | EST. CONC. | Q |
|---------|------------------|----|------------|---|
| 1. | No tics detected | | | |

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SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

Lab Name: GCAL Sample ID: SKGWFB1009
 Lab Code: LA024 Case No.: _____ Contract: _____
 SAS No.: _____ SDG No.: 204031909 Lab File ID: 2040326/S5029
 Matrix: Water Lab Sample ID: 20403190908
 Sample wt/vol: 1000 Units: mL Date Collected: 03/18/04 Time: 1702
 Level: (low/med) _____ Date Received: 03/19/04
 % Moisture: _____ decanted: (Y/N) _____ Date Extracted: 3/22/04
 GC Column: DB-5MS-30M ID: .25 (mm) Date Analyzed: 03/26/04 Time: 1855
 Concentrated Extract Volume: 1000 (µL) Dilution Factor: 1 Analyst: JAR3
 Injection Volume: 1.0 (µL) Prep Method: _____
 GPC Cleanup: (Y/N) N pH: _____ Analytical Method: OLMO 4.2
 Instrument ID: MSSV2

CONCENTRATION UNITS: ug/LPrep Batch: 271164 Analytical Batch: 271781

| CAS NO. | COMPOUND | RESULT | Q | MDL | PQL |
|-----------|-----------------------------|--------|---|-------|------|
| 95-95-4 | 2,4,5-Trichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 88-06-2 | 2,4,6-Trichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 120-83-2 | 2,4-Dichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 51-28-5 | 2,4-Dinitrophenol | 25.0 | U | 0.010 | 25.0 |
| 121-14-2 | 2,4-Dinitrotoluene | 10.0 | U | 0.010 | 10.0 |
| 606-20-2 | 2,6-Dinitrotoluene | 10.0 | U | 0.010 | 10.0 |
| 91-58-7 | 2-Chloronaphthalene | 10.0 | U | 0.010 | 10.0 |
| 95-57-8 | 2-Chlorophenol | 10.0 | U | 0.010 | 10.0 |
| 91-57-6 | 2-Methylnaphthalene | 10.0 | U | 0.010 | 10.0 |
| 88-74-4 | 2-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 88-75-5 | 2-Nitrophenol | 10.0 | U | 0.010 | 10.0 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 10.0 | U | 0.010 | 10.0 |
| 99-09-2 | 3-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 534-52-1 | 2-Methyl-4,6-dinitrophenol | 25.0 | U | 0.010 | 25.0 |
| 59-50-7 | 4-Chloro-3-methylphenol | 10.0 | U | 0.010 | 10.0 |
| 106-47-8 | 4-Chloroaniline | 10.0 | U | 0.010 | 10.0 |
| 7005-72-3 | 4-Chlorophenyl-phenylether | 10.0 | U | 0.010 | 10.0 |
| 106-44-5 | 4-Methylphenol (p-Cresol) | 10.0 | U | 0.010 | 10.0 |
| 83-32-9 | Acenaphthene | 10.0 | U | 0.010 | 10.0 |
| 208-96-8 | Acenaphthylene | 10.0 | U | 0.010 | 10.0 |
| 120-12-7 | Anthracene | 10.0 | U | 0.010 | 10.0 |
| 56-55-3 | Benzo(a)anthracene | 10.0 | U | 0.010 | 10.0 |
| 50-32-8 | Benzo(a)pyrene | 10.0 | U | 0.010 | 10.0 |
| 205-99-2 | Benzo(b)fluoranthene | 10.0 | U | 0.010 | 10.0 |
| 191-24-2 | Benzo(g,h,i)perylene | 10.0 | U | 0.010 | 10.0 |
| 207-08-9 | Benzo(k)fluoranthene | 10.0 | U | 0.010 | 10.0 |
| 111-91-1 | Bis(2-Chloroethoxy)methane | 10.0 | U | 0.010 | 10.0 |
| 111-44-4 | Bis(2-Chloroethyl)ether | 10.0 | U | 0.010 | 10.0 |
| 108-60-1 | bis(2-Chloroisopropyl)ether | 10.0 | U | 0.010 | 10.0 |
| 117-81-7 | bis(2-ethylhexyl)phthalate | 10.0 | U | 0.010 | 10.0 |

SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

| | |
|---|---|
| Lab Name: <u>GCAL</u> | Sample ID: <u>SKGWFB1009</u> |
| Lab Code: <u>LA024</u> Case No.: _____ | Contract: _____ |
| SAS No.: _____ SDG No.: <u>204031909</u> | Lab File ID: <u>2040326/S5029</u> |
| Matrix: <u>Water</u> | Lab Sample ID: <u>20403190908</u> |
| Sample wt/Vol: <u>1000</u> Units: <u>mL</u> | Date Collected: <u>03/18/04</u> Time: <u>1702</u> |
| Level: (low/med) _____ | Date Received: <u>03/19/04</u> |
| % Moisture: _____ Decanted: (Y/N) _____ | Date Extracted: <u>3/22/04</u> |
| GC Column: <u>DB-5MS-30M</u> ID: <u>25</u> (mm) | Date Analyzed: <u>03/26/04</u> Time: <u>1855</u> |
| Concentrated Extract Volume: <u>1000</u> (µL) | Dilution Factor: <u>1</u> Analyst: <u>JAR3</u> |
| Injection Volume: <u>1.0</u> (µL) | Prep Method: _____ |
| GPC Cleanup: (Y/N) <u>N</u> pH: _____ | Analytical Method: <u>OLMO 4.2</u> |
| | Instrument ID: <u>MSSV2</u> |

CONCENTRATION UNITS: µg/L

Prep Batch: 271164 Analytical Batch: 271781

| CAS NO. | COMPOUND | RESULT | Q | MDL | PQL |
|----------|----------------------------|--------|---|-------|------|
| 101-55-3 | 4-Bromophenyl-phenylether | 10.0 | U | 0.010 | 10.0 |
| 85-68-7 | Butylbenzylphthalate | 10.0 | U | 0.010 | 10.0 |
| 66-74-8 | Carbazole | 10.0 | U | 0.010 | 10.0 |
| 218-01-9 | Chrysene | 10.0 | U | 0.010 | 10.0 |
| 84-74-2 | Di-n-butylphthalate | 10.0 | U | 0.010 | 10.0 |
| 117-84-0 | Di-n-octylphthalate | 10.0 | U | 0.010 | 10.0 |
| 53-70-3 | Dibenz(a,h)anthracene | 10.0 | U | 0.010 | 10.0 |
| 132-64-9 | Dibenzofuran | 10.0 | U | 0.010 | 10.0 |
| 84-66-2 | Diethylphthalate | 10.0 | U | 0.010 | 10.0 |
| 131-11-3 | Dimethyl-phthalate | 10.0 | U | 0.010 | 10.0 |
| 105-67-9 | 2,4-Dimethylphenol | 10.0 | U | 0.010 | 10.0 |
| 206-44-0 | Fluoranthene | 10.0 | U | 0.010 | 10.0 |
| 86-73-7 | Fluorene | 10.0 | U | 0.010 | 10.0 |
| 118-74-1 | Hexachlorobenzene | 10.0 | U | 0.010 | 10.0 |
| 87-68-3 | Hexachlorobutadiene | 10.0 | U | 0.010 | 10.0 |
| 77-47-4 | Hexachlorocyclopentadiene | 10.0 | U | 0.010 | 10.0 |
| 67-72-1 | Hexachloroethane | 10.0 | U | 0.010 | 10.0 |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | 10.0 | U | 0.010 | 10.0 |
| 78-59-1 | Isophorone | 10.0 | U | 0.010 | 10.0 |
| 91-20-3 | Naphthalene | 10.0 | U | 0.010 | 10.0 |
| 100-01-6 | 4-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 98-95-3 | Nitrobenzene | 10.0 | U | 0.010 | 10.0 |
| 100-02-7 | 4-Nitrophenol | 25.0 | U | 0.010 | 25.0 |
| 87-86-5 | Pentachlorophenol | 25.0 | U | 0.010 | 25.0 |
| 85-01-8 | Phenanthrene | 10.0 | U | 0.010 | 10.0 |
| 108-95-2 | Phenol | 10.0 | U | 0.010 | 10.0 |
| 129-00-0 | Pyrene | 10.0 | U | 0.010 | 10.0 |
| 621-64-7 | N-Nitroso-di-n-propylamine | 10.0 | U | 0.010 | 10.0 |
| 86-30-6 | N-Nitrosodiphenylamine | 10.0 | U | 0.010 | 10.0 |
| 95-48-7 | o-Cresol | 10.0 | U | 0.010 | 10.0 |

1F
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

| | |
|--|---|
| Lab Name: <u>GCAL</u> | Sample ID: <u>SKGWFB1009</u> |
| Lab Code: <u>LA024</u> <u>2</u> Case No.: _____ | Contract: _____ |
| SAS No.: _____ SDG No.: <u>204031909</u> | Lab File ID: <u>S5029</u> |
| Matrix: <u>Water</u> | Lab Sample ID: <u>20403190908</u> |
| Sample wt/vol: _____ Units: _____ | Date Collected: <u>03/18/04</u> Time: <u>1702</u> |
| Level: (low/med) _____ | Date Received: <u>03/19/04</u> |
| % Moisture: not dec. _____ | Date Extracted: <u>3/22/04</u> |
| GC Column: <u>RTX-5MS-30</u> ID: <u>.53</u> (mm) | Date Analyzed: <u>03/26/04</u> Time: <u>1855</u> |
| Concentrated Extract Volume: <u>1000</u> (µL) | Dilution Factor: <u>1</u> Analyst: <u>RLW</u> |
| Injection Volume: <u>1.0</u> (µL) | Prep Method: _____ |
| GPC Cleanup: (Y/N) <u>N</u> pH: _____ | Analytical Method: <u>SW-846 8270C</u> |
| | Instrument ID: <u>MSSV2</u> |

Number TICs Found : 0

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | RT | EST. CONC. | Q |
|---------|------------------|----|------------|---|
| 1. | No tics detected | | | |

5/17/04
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SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

Lab Name: GCAL Sample ID: SKGW06R1009
 Lab Code: LA024 Case No.: _____ Contract: _____
 SAS No.: _____ SDG No.: 204031909 Lab File ID: 2040326/S5030
 Matrix: Water Lab Sample ID: 20403220801
 Sample wt/vol: 1000 Units: mL Date Collected: 03/16/04 Time: 1458
 Level: (low/med) _____ Date Received: 03/19/04
 % Moisture: _____ decanted: (Y/N) _____ Date Extracted: 3/22/04
 GC Column: DB-5MS-30M ID: 25 (mm) Date Analyzed: 03/26/04 Time: 1922
 Concentrated Extract Volume: 1000 (µL) Dilution Factor: 1 Analyst: JAR3
 Injection Volume: 1.0 (µL) Prep Method: _____
 GPC Cleanup: (Y/N) N pH: _____ Analytical Method: OLMO 4.2
 Instrument ID: MSSV2

CONCENTRATION UNITS: µg/L

Prep Batch: 271164 Analytical Batch: 271781

| CAS NO. | COMPOUND | RESULT | Q | MDL | PQL |
|-----------|-----------------------------|--------|---|-------|------|
| 95-95-4 | 2,4,5-Trichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 88-06-2 | 2,4,6-Trichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 120-83-2 | 2,4-Dichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 51-28-5 | 2,4-Dinitrophenol | 25.0 | U | 0.010 | 25.0 |
| 121-14-2 | 2,4-Dinitrotoluene | 10.0 | U | 0.010 | 10.0 |
| 606-20-2 | 2,6-Dinitrotoluene | 10.0 | U | 0.010 | 10.0 |
| 91-58-7 | 2-Chloronaphthalene | 10.0 | U | 0.010 | 10.0 |
| 95-57-8 | 2-Chlorophenol | 10.0 | U | 0.010 | 10.0 |
| 91-57-6 | 2-Methylnaphthalene | 10.0 | U | 0.010 | 10.0 |
| 88-74-4 | 2-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 88-75-5 | 2-Nitrophenol | 10.0 | U | 0.010 | 10.0 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 10.0 | U | 0.010 | 10.0 |
| 99-09-2 | 3-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 534-52-1 | 2-Methyl-4,6-dinitrophenol | 25.0 | U | 0.010 | 25.0 |
| 59-50-7 | 4-Chloro-3-methylphenol | 10.0 | U | 0.010 | 10.0 |
| 106-47-8 | 4-Chloroaniline | 10.0 | U | 0.010 | 10.0 |
| 7005-72-3 | 4-Chlorophenyl-phenylether | 10.0 | U | 0.010 | 10.0 |
| 106-44-5 | 4-Methylphenol p-Cresol | 10.0 | U | 0.010 | 10.0 |
| 83-32-9 | Acenaphthene | 10.0 | U | 0.010 | 10.0 |
| 208-96-8 | Acenaphthylene | 10.0 | U | 0.010 | 10.0 |
| 120-12-7 | Anthracene | 10.0 | U | 0.010 | 10.0 |
| 56-55-3 | Benzo(a)anthracene | 10.0 | U | 0.010 | 10.0 |
| 50-32-8 | Benzo(a)pyrene | 10.0 | U | 0.010 | 10.0 |
| 205-99-2 | Benzo(b)fluoranthene | 10.0 | U | 0.010 | 10.0 |
| 191-24-2 | Benzo(g,h)perylene | 10.0 | U | 0.010 | 10.0 |
| 207-08-9 | Benzo(k)fluoranthene | 10.0 | U | 0.010 | 10.0 |
| 111-91-1 | Bis(2-Chloroethoxy)methane | 10.0 | U | 0.010 | 10.0 |
| 111-44-4 | Bis(2-Chloroethyl)ether | 10.0 | U | 0.010 | 10.0 |
| 108-60-1 | bis(2-Chloroisopropyl)ether | 10.0 | U | 0.010 | 10.0 |
| 117-81-7 | bis(2-ethylhexyl)phthalate | 10.0 | U | 0.010 | 10.0 |

SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

Lab Name: GCAL Sample ID: SKGW06R1009
 Lab Code: LA024 Case No.: _____ Contract: _____
 SAS No.: _____ SDG No.: 204031909 Lab File ID: 2040326/S5030
 Matrix: Water Lab Sample ID: 20403220801
 Sample wt/vol: 1000 Units: mL Date Collected: 03/16/04 Time: 1458
 Level: (low/med) _____ Date Received: 03/19/04
 % Moisture: _____ decanted: (Y/N) _____ Date Extracted: 3/22/04
 GC Column: DB-5MS-30M ID: .25 (mm) Date Analyzed: 03/26/04 Time: 1922
 Concentrated Extract Volume: 1000 (μL) Dilution Factor: 1 Analyst: JAR3
 Injection Volume: 1.0 (μL) Prep Method: _____
 GPC Cleanup: (Y/N) N pH: _____ Analytical Method: OLMO 4.2
 Instrument ID: MSSV2
 CONCENTRATION UNITS: ug/L Prep Batch: 271164 Analytical Batch: 271781

| CAS NO. | COMPOUND | RESULT | Q | MDL | PQL |
|----------|----------------------------|--------|---|-------|------|
| 101-55-3 | 4-Bromophenyl-phenylether | 10.0 | U | 0.010 | 10.0 |
| 85-68-7 | Butylbenzylphthalate | 10.0 | U | 0.010 | 10.0 |
| 86-74-8 | Carbazole | 10.0 | U | 0.010 | 10.0 |
| 218-01-9 | Chrysene | 10.0 | U | 0.010 | 10.0 |
| 84-74-2 | Di-n-butylphthalate | 10.0 | U | 0.010 | 10.0 |
| 117-84-0 | Di-n-octylphthalate | 10.0 | U | 0.010 | 10.0 |
| 53-70-3 | Dibenz(a,h)anthracene | 10.0 | U | 0.010 | 10.0 |
| 132-64-9 | Dibenzofuran | 10.0 | U | 0.010 | 10.0 |
| 84-66-2 | Diethylphthalate | 10.0 | U | 0.010 | 10.0 |
| 131-11-3 | Dimethyl-phthalate | 10.0 | U | 0.010 | 10.0 |
| 105-67-9 | 2,4-Dimethylphenol | 10.0 | U | 0.010 | 10.0 |
| 206-44-0 | Fluoranthene | 10.0 | U | 0.010 | 10.0 |
| 86-73-7 | Fluorene | 10.0 | U | 0.010 | 10.0 |
| 118-74-1 | Hexachlorobenzene | 10.0 | U | 0.010 | 10.0 |
| 87-68-3 | Hexachlorobutadiene | 10.0 | U | 0.010 | 10.0 |
| 77-47-4 | Hexachlorocyclopentadiene | 10.0 | U | 0.010 | 10.0 |
| 67-72-1 | Hexachloroethane | 10.0 | U | 0.010 | 10.0 |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | 10.0 | U | 0.010 | 10.0 |
| 78-59-1 | Isophorone | 10.0 | U | 0.010 | 10.0 |
| 91-20-3 | Naphthalene | 10.0 | U | 0.010 | 10.0 |
| 100-01-6 | 4-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 98-95-3 | Nitrobenzene | 10.0 | U | 0.010 | 10.0 |
| 100-02-7 | 4-Nitrophenol | 25.0 | U | 0.010 | 25.0 |
| 87-86-5 | Pentachlorophenol | 25.0 | U | 0.010 | 25.0 |
| 85-01-8 | Phenanthrene | 10.0 | U | 0.010 | 10.0 |
| 108-95-2 | Phenol | 10.0 | U | 0.010 | 10.0 |
| 129-00-0 | Pyrene | 10.0 | U | 0.010 | 10.0 |
| 621-64-7 | N-Nitroso-di-n-propylamine | 10.0 | U | 0.010 | 10.0 |
| 86-30-6 | N-Nitrosodiphenylamine | 10.0 | U | 0.010 | 10.0 |
| 95-48-7 | o-Cresol | 10.0 | U | 0.010 | 10.0 |

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SEMI-VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

| | |
|--|---|
| Lab Name: <u>GCAL</u> | Sample ID: <u>SKGW06R1009</u> |
| Lab Code: <u>LA024</u> <u>2</u> Case No.: _____ | Contract: _____ |
| SAS No.: _____ SDG No.: <u>204031909</u> | Lab File ID: <u>S5030</u> |
| Matrix: <u>Water</u> | Lab Sample ID: <u>20403220801</u> |
| Sample wt/vol: _____ Units _____ | Date Collected: <u>03/16/04</u> Time: <u>1458</u> |
| Level: (low/med) _____ | Date Received: <u>03/19/04</u> |
| % Moisture: not dec. _____ | Date Extracted: <u>3/22/04</u> |
| GC Column: <u>RTX-SMS-30</u> ID: <u>.53</u> (mm) | Date Analyzed: <u>03/26/04</u> Time: <u>1922</u> |
| Concentrated Extract Volume: <u>*000</u> (µL) | Dilution Factor: <u>1</u> Analyst: <u>RLW</u> |
| Injection Volume: <u>*0</u> (µL) | Prep Method: _____ |
| GPC Cleanup: (Y/N) <u>N</u> pH: _____ | Analytical Method: <u>SW-846 8270C</u> |
| | Instrument ID: <u>MSSV2</u> |

Number TICs Found: 0

CONCENTRATION UNITS

| CAS NO. | COMPOUND | RT | EST. CONC. | Q |
|---------|------------------|----|------------|---|
| 1. | No pics detected | | | |

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SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

Lab Name: GCAL Sample ID: SKGW07R1009
 Lab Code: LA024 Case No.: _____ Contract: _____
 SAS No.: _____ SDG No.: 204031909 Lab File ID: 2040326/S5031
 Matrix: Water Lab Sample ID: 20403220802
 Sample wt/vol: 1000 Units: mL Date Collected: 03/16/04 Time: 1552
 Level: (low/med) _____ Date Received: 03/19/04
 % Moisture: _____ decanted: (Y/N) _____ Date Extracted: 3/22/04
 GC Column: DB-5MS-30M ID: .25 (mm) Date Analyzed: 03/26/04 Time: 1950
 Concentrated Extract Volume: 1000 (µL) Dilution Factor: 1 Analyst: JAR3
 Injection Volume: 1.0 (µL) Prep Method: _____
 GPC Cleanup: (Y/N) N pH: _____ Analytical Method: OLMO 4.2
 Instrument ID: MSSV2

CONCENTRATION UNITS: ug/LPrep Batch: 271164 Analytical Batch: 271781

| CAS NO. | COMPOUND | RESULT | Q | MDL | PQL |
|-----------|-----------------------------|--------|---|-------|------|
| 95-95-4 | 2,4,5-Trichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 88-06-2 | 2,4,6-Trichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 120-83-2 | 2,4-Dichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 51-28-5 | 2,4-Dinitrophenol | 25.0 | U | 0.010 | 25.0 |
| 121-14-2 | 2,4-Dinitrotoluene | 10.0 | U | 0.010 | 10.0 |
| 606-20-2 | 2,6-Dinitrotoluene | 10.0 | U | 0.010 | 10.0 |
| 91-58-7 | 2-Chloronaphthalene | 10.0 | U | 0.010 | 10.0 |
| 95-57-8 | 2-Chlorophenol | 10.0 | U | 0.010 | 10.0 |
| 91-57-6 | 2-Methylnaphthalene | 10.0 | U | 0.010 | 10.0 |
| 88-74-4 | 2-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 88-75-5 | 2-Nitrophenol | 10.0 | U | 0.010 | 10.0 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 10.0 | U | 0.010 | 10.0 |
| 99-09-2 | 3-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 534-52-1 | 2-Methyl-4,6-dinitrophenol | 25.0 | U | 0.010 | 25.0 |
| 59-50-7 | 4-Chloro-3-methylphenol | 10.0 | U | 0.010 | 10.0 |
| 106-47-8 | 4-Chloroaniline | 10.0 | U | 0.010 | 10.0 |
| 7005-72-3 | 4-Chlorophenyl-phenylether | 10.0 | U | 0.010 | 10.0 |
| 106-44-5 | 4-Methylphenol (p-Cresol) | 10.0 | U | 0.010 | 10.0 |
| 83-32-9 | Acenaphthene | 10.0 | U | 0.010 | 10.0 |
| 208-96-8 | Acenaphthylene | 10.0 | U | 0.010 | 10.0 |
| 120-12-7 | Anthracene | 10.0 | U | 0.010 | 10.0 |
| 56-55-3 | Benzo(a)anthracene | 10.0 | U | 0.010 | 10.0 |
| 50-32-8 | Benzo(a)pyrene | 10.0 | U | 0.010 | 10.0 |
| 205-99-2 | Benzo(b)fluoranthene | 10.0 | U | 0.010 | 10.0 |
| 191-24-2 | Benzo(g,h,i)perylene | 10.0 | U | 0.010 | 10.0 |
| 207-08-9 | Benzo(k)fluoranthene | 10.0 | U | 0.010 | 10.0 |
| 111-91-1 | Bis(2-Chloroethoxy)methane | 10.0 | U | 0.010 | 10.0 |
| 111-44-4 | Bis(2-Chloroethyl)ether | 10.0 | U | 0.010 | 10.0 |
| 108-60-1 | bis(2-Chloroisopropyl)ether | 10.0 | U | 0.010 | 10.0 |
| 117-81-7 | bis(2-ethylhexyl)phthalate | 10.0 | U | 0.010 | 10.0 |

SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

Lab Name: GCAL Sample ID: SKGW07R1009

Lab Code: LA024 Case No.: _____ Contract: _____

SAS No.: _____ SDG No.: 204031909 Lab File ID: 2040326/S5031

Matrix: Water Lab Sample ID: 20403220802

Sample wt/vol: 1000 Units: mL Date Collected: 03/16/04 Time: 1552

Level: (low/med) _____ Date Received: 03/19/04

% Moisture: _____ decanted: (Y/N) _____ Date Extracted: 3/22/04

GC Column: DB-5MS-30M ID: 25 (mm) Date Analyzed: 03/26/04 Time: 1950

Concentrated Extract Volume: 1000 (µL) Dilution Factor: 1 Analyst: JAR3

Injection Volume: 10 (µL) Prep Method: _____

GPC Cleanup: (Y/N) N pH: _____ Analytical Method: OLMO 4.2

Instrument ID: MSSV2

Prep Batch: 271164 Analytical Batch: 271781

CONCENTRATION UNITS: µg/L

| CAS NO. | COMPOUND | RESULT | Q | MDL | PQL |
|----------|----------------------------|--------|---|-------|------|
| 101-55-3 | 4-Bromophenyl-phenylether | 10.0 | U | 0.010 | 10.0 |
| 85-68-7 | Butylbenzylphthalate | 10.0 | U | 0.010 | 10.0 |
| 86-74-8 | Carbazole | 10.0 | U | 0.010 | 10.0 |
| 218-01-9 | Chrysene | 10.0 | U | 0.010 | 10.0 |
| 84-74-2 | Di-n-butylphthalate | 10.0 | U | 0.010 | 10.0 |
| 117-84-0 | Di-n-octylphthalate | 10.0 | U | 0.010 | 10.0 |
| 53-70-3 | Dibenz(a,h)anthracene | 10.0 | U | 0.010 | 10.0 |
| 132-64-9 | Dibenzofuran | 10.0 | U | 0.010 | 10.0 |
| 84-66-2 | Diethylphthalate | 10.0 | U | 0.010 | 10.0 |
| 131-11-3 | Dimethyl-phthalate | 10.0 | U | 0.010 | 10.0 |
| 105-67-9 | 2,4-Dimethylphenol | 10.0 | U | 0.010 | 10.0 |
| 206-44-0 | Fluoranthene | 10.0 | U | 0.010 | 10.0 |
| 86-73-7 | Fluorene | 10.0 | U | 0.010 | 10.0 |
| 118-74-1 | Hexachlorobenzene | 10.0 | U | 0.010 | 10.0 |
| 87-68-3 | Hexachlorocycloadiene | 10.0 | U | 0.010 | 10.0 |
| 77-47-4 | Hexachlorocyclopentadiene | 10.0 | U | 0.010 | 10.0 |
| 87-72-1 | Hexachloroethane | 10.0 | U | 0.010 | 10.0 |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | 10.0 | U | 0.010 | 10.0 |
| 78-59-1 | Isophorone | 10.0 | U | 0.010 | 10.0 |
| 91-20-3 | Naphthalene | 10.0 | U | 0.010 | 10.0 |
| 100-01-6 | 4-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 98-95-3 | Nitrobenzene | 10.0 | U | 0.010 | 10.0 |
| 100-02-7 | 4-Nitrophenol | 25.0 | U | 0.010 | 25.0 |
| 87-86-5 | Pentachlorophenol | 25.0 | U | 0.010 | 25.0 |
| 85-01-8 | Phenanthrene | 10.0 | U | 0.010 | 10.0 |
| 108-95-2 | Phenol | 10.0 | U | 0.010 | 10.0 |
| 129-00-0 | Pyrene | 10.0 | U | 0.010 | 10.0 |
| 621-64-7 | N-Nitroso-di-n-propylamine | 10.0 | U | 0.010 | 10.0 |
| 86-30-6 | N-Nitrosodiphenylamine | 10.0 | U | 0.010 | 10.0 |
| 95-48-7 | o-Cresol | 10.0 | U | 0.010 | 10.0 |

SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

Lab Name: GCAL Sample ID: SKGW07R1009
Lab Code: LA024 2 Case No.: _____ Contract: _____
SAS No.: _____ SDG No.: 204031909 Lab File ID: S5031
Matrix: Water Lab Sample ID: 20403220802
Sample wt/vol: _____ Units: _____ Date Collected: 03/16/04 Time: 1552
Level: (low/med) _____ Date Received: 03/19/04
% Moisture: not dec. _____ Date Extracted: 3/22/04
GC Column: RTX-5MS-30 ID: .53 (mm) Date Analyzed: 03/26/04 Time: 1950
Concentrated Extract Volume: 1000 (µL) Dilution Factor: 1 Analyst: RLW
Injection Volume: 1.0 (µL) Prep Method: _____
GPC Cleanup: (Y/N) N pH: _____ Analytical Method: SW-846 8270C
Instrument ID: MSSV2

Number TICs Found : 0

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | RT | EST. CONC. | Q |
|---------|------------------|----|------------|---|
| 1. | No tics detected | | | |

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SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

Lab Name: GCAL Sample ID: SKGW591009
 Lab Code: LA024 Case No.: _____ Contract: _____
 SAS No.: _____ SDG No.: 204031909 Lab File ID: 2040326/S5032
 Matrix: Water Lab Sample ID: 20403220803
 Sample wt/vol: 1000 Units: mL Date Collected: 03/17/04 Time: 1050
 Level: (low/med) _____ Date Received: 03/19/04
 % Moisture: _____ decanted: (Y/N) _____ Date Extracted: 3/22/04
 GC Column: DB-5MS-30M ID: 25 (mm) Date Analyzed: 03/26/04 Time: 2018
 Concentrated Extract Volume 1000 (µL) Dilution Factor: 1 Analyst: JAR3
 Injection Volume: 1.0 (µL) Prep Method: _____
 GPC Cleanup: (Y/N) N pH: _____ Analytical Method: OLMO 4.2
 Instrument ID: MSSV2

CONCENTRATION UNITS µg/L

Prep Batch: 271164 Analytical Batch: 271781

| CAS NO. | COMPOUND | RESULT | Q | MDL | PQL |
|-----------|-----------------------------|--------|---|-------|------|
| 95-95-4 | 2,4,5-Trichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 88-06-2 | 2,4,6-Trichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 120-83-2 | 2,4-Dichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 51-28-5 | 2,4-Dinitrophenol | 25.0 | U | 0.010 | 25.0 |
| 121-14-2 | 2,4-Dinitrobenzene | 10.0 | U | 0.010 | 10.0 |
| 606-20-2 | 2,6-Dinitrobenzene | 10.0 | U | 0.010 | 10.0 |
| 91-58-7 | 2-Chloronaphthalene | 10.0 | U | 0.010 | 10.0 |
| 95-57-8 | 2-Chlorophenol | 10.0 | U | 0.010 | 10.0 |
| 91-57-6 | 2-Methylnaphthalene | 10.0 | U | 0.010 | 10.0 |
| 88-74-4 | 2-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 88-75-5 | 2-Nitrophenol | 10.0 | U | 0.010 | 10.0 |
| 91-04-1 | 3,3'-Dichlorobenzidine | 10.0 | U | 0.010 | 10.0 |
| 99-08-2 | 3-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 534-52-1 | 2-Methyl-4-nitrophenol | 25.0 | U | 0.010 | 25.0 |
| 58-50-7 | 4-Chloro-3-methylphenol | 10.0 | U | 0.010 | 10.0 |
| 106-47-8 | 4-Chloroaniline | 10.0 | U | 0.010 | 10.0 |
| 7005-72-3 | 4-Chlorophenyl-phenylether | 10.0 | U | 0.010 | 10.0 |
| 106-44-5 | 4-Methylphenol (p-Cresol) | 10.0 | U | 0.010 | 10.0 |
| 83-32-9 | Acenaphthene | 10.0 | U | 0.010 | 10.0 |
| 208-96-8 | Acenaphthylene | 10.0 | U | 0.010 | 10.0 |
| 120-12-7 | Anthracene | 10.0 | U | 0.010 | 10.0 |
| 56-55-3 | Benzo(a)anthracene | 10.0 | U | 0.010 | 10.0 |
| 50-32-8 | Benzo(a)pyrene | 10.0 | U | 0.010 | 10.0 |
| 205-99-2 | Benzo(b)fluoranthene | 10.0 | U | 0.010 | 10.0 |
| 191-24-2 | Benzo(g,h)perylene | 10.0 | U | 0.010 | 10.0 |
| 207-08-9 | Benzo(k)fluoranthene | 10.0 | U | 0.010 | 10.0 |
| 111-91-1 | Bis(2-Chloroethoxy)methane | 10.0 | U | 0.010 | 10.0 |
| 111-44-4 | Bis(2-Chloroethyl)ether | 10.0 | U | 0.010 | 10.0 |
| 108-60-1 | bis(2-Chloroisopropyl)ether | 10.0 | U | 0.010 | 10.0 |
| 117-81-7 | bis(2-ethoxyethyl)phthalate | 10.0 | U | 0.010 | 10.0 |

SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

Lab Name: GCAL Sample ID: SKGW591009

Lab Code: LA024 Case No.: _____ Contract: _____

SAS No.: _____ SDG No.: 204031909 Lab File ID: 2040326/S5032

Matrix: Water Lab Sample ID: 20403220803

Sample wt/vol: 1000 Units: mL Date Collected: 03/17/04 Time: 1050

Level: (low/med) _____ Date Received: 03/19/04

% Moisture: _____ decanted: (Y/N) _____ Date Extracted: 3/22/04

GC Column: DB-5MS-30M ID: .25 (mm) Date Analyzed: 03/26/04 Time: 2018

Concentrated Extract Volume: 1000 (µL) Dilution Factor: 1 Analyst: JAR3

Injection Volume: 1.0 (µL) Prep Method: _____

GPC Cleanup: (Y/N) N pH: _____ Analytical Method: OLMO 4.2

Instrument ID: MSSV2

Prep Batch: 271164 Analytical Batch: 271781

CONCENTRATION UNITS: ug/L

| CAS NO. | COMPOUND | RESULT | Q | MDL | PQL |
|----------|----------------------------|--------|---|-------|------|
| 101-55-3 | 4-Bromophenyl-phenylether | 10.0 | U | 0.010 | 10.0 |
| 85-68-7 | Butylbenzylphthalate | 10.0 | U | 0.010 | 10.0 |
| 86-74-8 | Carbazole | 10.0 | U | 0.010 | 10.0 |
| 218-01-9 | Chrysene | 10.0 | U | 0.010 | 10.0 |
| 84-74-2 | Di-n-butylphthalate | 10.0 | U | 0.010 | 10.0 |
| 117-84-0 | Di-n-octylphthalate | 10.0 | U | 0.010 | 10.0 |
| 53-70-3 | Dibenz(a,h)anthracene | 10.0 | U | 0.010 | 10.0 |
| 132-64-9 | Dibenzofuran | 10.0 | U | 0.010 | 10.0 |
| 84-66-2 | Diethylphthalate | 10.0 | U | 0.010 | 10.0 |
| 131-11-3 | Dimethyl-phthalate | 10.0 | U | 0.010 | 10.0 |
| 105-67-9 | 2,4-Dimethylphenol | 10.0 | U | 0.010 | 10.0 |
| 206-44-0 | Fluoranthene | 10.0 | U | 0.010 | 10.0 |
| 86-73-7 | Fluorene | 10.0 | U | 0.010 | 10.0 |
| 118-74-1 | Hexachlorobenzene | 10.0 | U | 0.010 | 10.0 |
| 87-68-3 | Hexachlorobutadiene | 10.0 | U | 0.010 | 10.0 |
| 77-47-4 | Hexachlorocyclopentadiene | 10.0 | U | 0.010 | 10.0 |
| 67-72-1 | Hexachloroethane | 10.0 | U | 0.010 | 10.0 |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | 10.0 | U | 0.010 | 10.0 |
| 78-59-1 | Isophorone | 10.0 | U | 0.010 | 10.0 |
| 91-20-3 | Naphthalene | 10.0 | U | 0.010 | 10.0 |
| 100-01-6 | 4-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 98-95-3 | Nitrobenzene | 10.0 | U | 0.010 | 10.0 |
| 100-02-7 | 4-Nitrophenol | 25.0 | U | 0.010 | 25.0 |
| 87-86-5 | Pentachlorophenol | 25.0 | U | 0.010 | 25.0 |
| 85-01-8 | Phenanthrene | 10.0 | U | 0.010 | 10.0 |
| 108-95-2 | Phenol | 10.0 | U | 0.010 | 10.0 |
| 129-00-0 | Pyrene | 10.0 | U | 0.010 | 10.0 |
| 621-64-7 | N-Nitroso-di-n-propylamine | 10.0 | U | 0.010 | 10.0 |
| 85-30-6 | N-Nitrosodiphenylamine | 10.0 | U | 0.010 | 10.0 |
| 95-48-7 | o-Cresol | 10.0 | U | 0.010 | 10.0 |

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435/13/09
msa

SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

| | |
|---|---|
| Lab Name: <u>GCAL</u> | Sample ID: <u>SKGW591009</u> |
| Lab Code: <u>LA024</u> <u>2</u> Case No.: _____ | Contract: _____ |
| SAS No.: _____ SDG No.: <u>204031909</u> | Lab File ID: <u>S5032</u> |
| Matrix: <u>Water</u> | Lab Sample ID: <u>20403220803</u> |
| Sample wt/vol: _____ Units: _____ | Date Collected: <u>03/17/04</u> Time: <u>1050</u> |
| Level: (low/med) _____ | Date Received: <u>03/19/04</u> |
| % Moisture: not dec. _____ | Date Extracted: <u>3/22/04</u> |
| GC Column: <u>RTX-5MS-30</u> ID: <u>53</u> (mm) | Date Analyzed: <u>03/26/04</u> Time: <u>2018</u> |
| Concentrated Extract Volume: <u>1000</u> (µL) | Dilution Factor: <u>1</u> Analyst: <u>RLW</u> |
| Injection Volume: <u>1.0</u> (µL) | Prep Method: _____ |
| GPC Cleanup: (Y/N) <u>N</u> pH: _____ | Analytical Method: <u>SW-846 8270C</u> |
| | Instrument ID: <u>MSSV2</u> |

Number TICs Found : 0

CONCENTRATION UNITS

| CAS NO. | COMPOUND | RT | EST. CONC. | Q |
|---------|------------------|----|------------|---|
| 1. | No TICs detected | | | |

5/17/04
RLW

SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

Lab Name: GCAL Sample ID: SKGW601009
 Lab Code: LA024 Case No.: _____ Contract: _____
 SAS No.: _____ SDG No.: 204031909 Lab File ID: 2040326/S5033
 Matrix: Water Lab Sample ID: 20403220804
 Sample wt/vol: 1000 Units: mL Date Collected: 03/17/04 Time: 1115
 Level: (low/med) _____ Date Received: 03/19/04
 % Moisture: _____ decanted: (Y/N) _____ Date Extracted: 3/22/04
 GC Column: DB-5MS-30M ID: .25 (mm) Date Analyzed: 03/26/04 Time: 2045
 Concentrated Extract Volume: 1000 (μL) Dilution Factor: 1 Analyst: JAR3
 Injection Volume: 1.0 (μL) Prep Method: _____
 GPC Cleanup: (Y/N) N pH: _____ Analytical Method: OLMO 4.2
 Instrument ID: MSSV2
 Prep Batch: 271164 Analytical Batch: 271781

CONCENTRATION UNITS: ug/L

| CAS NO. | COMPOUND | RESULT | Q | MDL | PQL |
|-----------|-----------------------------|--------|---|-------|------|
| 95-95-4 | 2,4,5-Trichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 88-06-2 | 2,4,6-Trichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 120-83-2 | 2,4-Dichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 51-28-5 | 2,4-Dinitrophenol | 25.0 | U | 0.010 | 25.0 |
| 121-14-2 | 2,4-Dinitrotoluene | 10.0 | U | 0.010 | 10.0 |
| 606-20-2 | 2,6-Dinitrotoluene | 10.0 | U | 0.010 | 10.0 |
| 91-58-7 | 2-Chloronaphthalene | 10.0 | U | 0.010 | 10.0 |
| 95-57-8 | 2-Chlorophenol | 10.0 | U | 0.010 | 10.0 |
| 91-57-6 | 2-Methylnaphthalene | 10.0 | U | 0.010 | 10.0 |
| 88-74-4 | 2-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 88-75-5 | 2-Nitrophenol | 10.0 | U | 0.010 | 10.0 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 10.0 | U | 0.010 | 10.0 |
| 99-09-2 | 3-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 534-52-1 | 2-Methyl-4,6-dinitrophenol | 25.0 | U | 0.010 | 25.0 |
| 59-50-7 | 4-Chloro-3-methylphenol | 10.0 | U | 0.010 | 10.0 |
| 106-47-8 | 4-Chloroaniline | 10.0 | U | 0.010 | 10.0 |
| 7005-72-3 | 4-Chlorophenyl-phenylether | 10.0 | U | 0.010 | 10.0 |
| 106-44-5 | 4-Methylphenol (p-Cresol) | 10.0 | U | 0.010 | 10.0 |
| 83-32-9 | Acenaphthene | 10.0 | U | 0.010 | 10.0 |
| 206-96-8 | Acenaphthylene | 10.0 | U | 0.010 | 10.0 |
| 120-12-7 | Anthracene | 10.0 | U | 0.010 | 10.0 |
| 56-55-3 | Benzo(a)anthracene | 10.0 | U | 0.010 | 10.0 |
| 50-32-8 | Benzo(a)pyrene | 10.0 | U | 0.010 | 10.0 |
| 205-99-2 | Benzo(b)fluoranthene | 10.0 | U | 0.010 | 10.0 |
| 191-24-2 | Benzo(g,h,i)perylene | 10.0 | U | 0.010 | 10.0 |
| 207-08-9 | Benzo(k)fluoranthene | 10.0 | U | 0.010 | 10.0 |
| 111-91-1 | Bis(2-Chloroethoxy)methane | 10.0 | U | 0.010 | 10.0 |
| 111-44-4 | Bis(2-Chloroethyl)ether | 10.0 | U | 0.010 | 10.0 |
| 108-60-1 | bis(2-Chloroisopropyl)ether | 10.0 | U | 0.010 | 10.0 |
| 117-81-7 | bis(2-ethylhexyl)phthalate | 10.0 | U | 0.010 | 10.0 |

5/17/07
msm

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SEMI-VOLATILE ORGANICS ANALYSIS DATA SHEET

Lab Name: GCAL Sample ID: SKGW601009
 Lab Code: LA024 Case No.: _____ Contract: _____
 SAS No.: _____ SCG No.: 204031909 Lab File ID: 2040326/S5033
 Matrix: Water Lab Sample ID: 20403220804
 Sample wt/vol: 1000 Units mL Date Collected: 03/17/04 Time: 1115
 Level: (low/med) _____ Date Received: 03/19/04
 % Moisture: _____ Decanted: (Y/N) _____ Date Extracted: 3/22/04
 GC Column: DB-5MS-30M ID: 25 (mm) Date Analyzed: 03/26/04 Time: 2045
 Concentrated Extract Volume 1000 (µL) Dilution Factor: 1 Analyst: JAR3
 Injection Volume: 10 (µL) Prep Method: _____
 GPC Cleanup: (Y/N) N pH: _____ Analytical Method: OLMO 4.2
 Instrument ID: MSSV2
 Prep Batch: 271164 Analytical Batch: 271781
 CONCENTRATION UNITS: µg/L

| CAS NO. | COMPOUND | RESULT | Q | MDL | PQL |
|----------|----------------------------|--------|---|-------|------|
| 101-55-3 | 4-Bromophenyl-phenylether | 10.0 | U | 0.010 | 10.0 |
| 85-68-7 | Butylbenzylphthalate | 10.0 | U | 0.010 | 10.0 |
| 86-74-8 | Carbazole | 10.0 | U | 0.010 | 10.0 |
| 218-01-8 | Chrysene | 10.0 | U | 0.010 | 10.0 |
| 84-74-2 | Di-n-butylphthalate | 10.0 | U | 0.010 | 10.0 |
| 117-84-0 | Di-n-octylphthalate | 10.0 | U | 0.010 | 10.0 |
| 53-70-3 | Dibenz(a,h)anthracene | 10.0 | U | 0.010 | 10.0 |
| 132-84-9 | Dibenzofuran | 10.0 | U | 0.010 | 10.0 |
| 84-66-2 | Diethylphthalate | 10.0 | U | 0.010 | 10.0 |
| 131-11-3 | Dimethyl-phthalate | 10.0 | U | 0.010 | 10.0 |
| 105-67-9 | 2,4-Dimethylphenol | 10.0 | U | 0.010 | 10.0 |
| 208-44-0 | Fluoranthene | 10.0 | U | 0.010 | 10.0 |
| 86-73-7 | Fluorene | 10.0 | U | 0.010 | 10.0 |
| 118-74-1 | Hexachlorobenzene | 10.0 | U | 0.010 | 10.0 |
| 87-68-3 | Hexachlorocyclopentadiene | 10.0 | U | 0.010 | 10.0 |
| 77-47-4 | Hexachlorocyclopentadiene | 10.0 | U | 0.010 | 10.0 |
| 67-72-1 | Hexachloroethane | 10.0 | U | 0.010 | 10.0 |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | 10.0 | U | 0.010 | 10.0 |
| 78-59-1 | Isophorone | 10.0 | U | 0.010 | 10.0 |
| 91-20-3 | Naphthalene | 10.0 | U | 0.010 | 10.0 |
| 100-01-6 | 4-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 98-95-3 | Nitrobenzene | 10.0 | U | 0.010 | 10.0 |
| 100-02-7 | 4-Nitrophenol | 25.0 | U | 0.010 | 25.0 |
| 87-86-5 | Pentachlorophenol | 25.0 | U | 0.010 | 25.0 |
| 85-01-8 | Phenanthrene | 10.0 | U | 0.010 | 10.0 |
| 108-95-2 | Phenol | 10.0 | U | 0.010 | 10.0 |
| 129-00-0 | Pyrene | 10.0 | U | 0.010 | 10.0 |
| 621-64-7 | N-Nitroso-Di-n-propylamine | 10.0 | U | 0.010 | 10.0 |
| 86-30-6 | N-Nitrosodiphenylamine | 10.0 | U | 0.010 | 10.0 |
| 95-48-7 | o-Cresol | 10.0 | U | 0.010 | 10.0 |

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SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

Lab Name: GCAL Sample ID: SKGW601009
Lab Code: LA024 2 Case No.: _____ Contract: _____
SAS No.: _____ SDG No.: 204031909 Lab File ID: S5033
Matrix: Water Lab Sample ID: 20403220804
Sample wt/vol: _____ Units: _____ Date Collected: 03/17/04 Time: 1115
Level: (low/med) _____ Date Received: 03/19/04
% Moisture: not dec. _____ Date Extracted: 3/22/04
GC Column: RTX-5MS-30 ID: .53 (mm) Date Analyzed: 03/26/04 Time: 2045
Concentrated Extract Volume: 1000 (µL) Dilution Factor: 1 Analyst: RLW
Injection Volume: 1.0 (µL) Prep Method: _____
GPC Cleanup: (Y/N) N pH: _____ Analytical Method: SW-846 8270C
Instrument ID: MSSV2

Number TICs Found : 0

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | RT | EST. CONC. | Q |
|---------|------------------|----|------------|---|
| 1. | No tics detected | | | |

5/17/04
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SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

| | |
|---|---|
| Lab Name: <u>GCAL</u> | Sample ID: <u>SKGW62A1009</u> |
| Lab Code: <u>LA024</u> Case No.: _____ | Contract: _____ |
| SAS No.: _____ SDG No.: <u>204031909</u> | Lab File ID: <u>2040326/S5034</u> |
| Matrix: <u>Water</u> | Lab Sample ID: <u>20403220805</u> |
| Sample wt/vol: <u>1000</u> Units: <u>mL</u> | Date Collected: <u>03/17/04</u> Time: <u>1150</u> |
| Level: (low/med) _____ | Date Received: <u>03/19/04</u> |
| % Moisture: _____ decanted: (Y/N) _____ | Date Extracted: <u>7/22/04</u> |
| GC Column: <u>DB-5MS-30M</u> ID: <u>25</u> (mm) | Date Analyzed: <u>03/26/04</u> Time: <u>2112</u> |
| Concentrated Extract Volume: <u>1000</u> (µL) | Dilution Factor: <u>1</u> Analyst: <u>JAR3</u> |
| Injection Volume: <u>1.0</u> (µL) | Prep Method: _____ |
| GPC Cleanup: (Y/N) <u>N</u> pH: _____ | Analytical Method: <u>OLMO 4.2</u> |
| CONCENTRATION UNITS: <u>µg/L</u> | Instrument ID: <u>MSSV2</u> |
| | Prep Batch: <u>271164</u> Analytical Batch: <u>271781</u> |

| CAS NO. | COMPOUND | RESULT | Q | MDL | PQL |
|-----------|-------------------------------|--------|---|-------|------|
| 95-95-4 | 2,4,5-Trichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 98-06-2 | 2,4,6-Trichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 120-83-2 | 2,4-Dichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 51-28-5 | 2,4-Dinitrophenol | 25.0 | U | 0.010 | 25.0 |
| 121-14-2 | 2,4-Dinitrotoluene | 10.0 | U | 0.010 | 10.0 |
| 806-20-2 | 2,6-Dinitrotoluene | 10.0 | U | 0.010 | 10.0 |
| 91-58-7 | 2-Chloronaphthalene | 10.0 | U | 0.010 | 10.0 |
| 95-57-8 | 2-Chlorophenol | 10.0 | U | 0.010 | 10.0 |
| 91-57-6 | 2-Methylnaphthalene | 10.0 | U | 0.010 | 10.0 |
| 88-74-4 | 2-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 88-75-5 | 2-Nitrophenol | 10.0 | U | 0.010 | 10.0 |
| 91-84-1 | 3,3'-Dichlorodiphenylmethane | 10.0 | U | 0.010 | 10.0 |
| 98-09-2 | 3-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 534-52-1 | 2-Methyl-4,6-dinitrophenol | 25.0 | U | 0.010 | 25.0 |
| 58-50-7 | 4-Chloro-3-methylphenol | 10.0 | U | 0.010 | 10.0 |
| 108-47-8 | 4-Chloroaniline | 10.0 | U | 0.010 | 10.0 |
| 7005-72-3 | 4-Chlorophenyl-phenylether | 10.0 | U | 0.010 | 10.0 |
| 108-44-5 | 4-Methylphenol (p-Cresol) | 10.0 | U | 0.010 | 10.0 |
| 83-32-9 | Acenaphthene | 10.0 | U | 0.010 | 10.0 |
| 208-96-8 | Acenaphthylene | 10.0 | U | 0.010 | 10.0 |
| 120-12-7 | Anthracene | 10.0 | U | 0.010 | 10.0 |
| 56-55-3 | Benzo(a)anthracene | 10.0 | U | 0.010 | 10.0 |
| 50-32-8 | Benzo(a)pyrene | 10.0 | U | 0.010 | 10.0 |
| 205-99-2 | Benzo(b)fluoranthene | 10.0 | U | 0.010 | 10.0 |
| 191-24-2 | Benzo(g,h,i)perylene | 10.0 | U | 0.010 | 10.0 |
| 207-08-9 | Benzo(k)fluoranthene | 10.0 | U | 0.010 | 10.0 |
| 111-91-1 | Bis(2-Chloroethoxy)methane | 10.0 | U | 0.010 | 10.0 |
| 111-44-4 | Bis(2-Chloroethyl)ether | 10.0 | U | 0.010 | 10.0 |
| 108-60-1 | bis(2-Chloroisopropyl)ether | 10.0 | U | 0.010 | 10.0 |
| 117-81-7 | bis(2-ethylhexyloxy)phthalate | 10.0 | U | 0.010 | 10.0 |

SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

Lab Name: GCAL Sample ID: SKGW62A1009
 Lab Code: LA024 Case No.: _____ Contract: _____
 SAS No.: _____ SDG No.: 204031909 Lab File ID: 2040326/S5034
 Matrix: Water Lab Sample ID: 20403220805
 Sample wt/vol: 1000 Units: mL Date Collected: 03/17/04 Time: 1150
 Level: (low/med) _____ Date Received: 03/19/04
 % Moisture: _____ decanted: (Y/N) _____ Date Extracted: 3/22/04
 GC Column: DB-5MS-30M ID: .25 (mm) Date Analyzed: 03/26/04 Time: 2112
 Concentrated Extract Volume: 1000 (µL) Dilution Factor: 1 Analyst: JAR3
 Injection Volume: 1.0 (µL) Prep Method: _____
 GPC Cleanup: (Y/N) N pH: _____ Analytical Method: OLMO 4.2
 Instrument ID: MSSV2
 Prep Batch: 271164 Analytical Batch: 271781

| CAS NO. | COMPOUND | RESULT | Q | MDL | PQL |
|----------|----------------------------|--------|---|-------|------|
| 101-55-3 | 4-Bromophenyl-phenylether | 10.0 | U | 0.010 | 10.0 |
| 85-68-7 | Butylbenzylphthalate | 10.0 | U | 0.010 | 10.0 |
| 86-74-8 | Carbazole | 10.0 | U | 0.010 | 10.0 |
| 218-01-9 | Chrysene | 10.0 | U | 0.010 | 10.0 |
| 84-74-2 | Di-n-butylphthalate | 10.0 | U | 0.010 | 10.0 |
| 117-84-0 | Di-n-octylphthalate | 10.0 | U | 0.010 | 10.0 |
| 53-70-3 | Dibenz(a,h)anthracene | 10.0 | U | 0.010 | 10.0 |
| 132-64-9 | Dibenzofuran | 10.0 | U | 0.010 | 10.0 |
| 84-66-2 | Diethylphthalate | 10.0 | U | 0.010 | 10.0 |
| 131-11-3 | Dimethyl-phthalate | 10.0 | U | 0.010 | 10.0 |
| 105-67-9 | 2,4-Dimethylphenol | 10.0 | U | 0.010 | 10.0 |
| 206-44-0 | Fluoranthene | 10.0 | U | 0.010 | 10.0 |
| 86-73-7 | Fluorene | 10.0 | U | 0.010 | 10.0 |
| 118-74-1 | Hexachlorobenzene | 10.0 | U | 0.010 | 10.0 |
| 87-68-3 | Hexachlorobutadiene | 10.0 | U | 0.010 | 10.0 |
| 77-47-4 | Hexachlorocyclopentadiene | 10.0 | U | 0.010 | 10.0 |
| 67-72-1 | Hexachloroethane | 10.0 | U | 0.010 | 10.0 |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | 10.0 | U | 0.010 | 10.0 |
| 78-59-1 | Isophorone | 10.0 | U | 0.010 | 10.0 |
| 91-20-3 | Naphthalene | 10.0 | U | 0.010 | 10.0 |
| 100-01-6 | 4-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 98-95-3 | Nitrobenzene | 10.0 | U | 0.010 | 10.0 |
| 100-02-7 | 4-Nitrophenol | 25.0 | U | 0.010 | 25.0 |
| 87-86-5 | Pentachlorophenol | 25.0 | U | 0.010 | 25.0 |
| 85-01-8 | Phenanthrene | 10.0 | U | 0.010 | 10.0 |
| 108-95-2 | Phenol | 10.0 | U | 0.010 | 10.0 |
| 129-00-0 | Pyrene | 10.0 | U | 0.010 | 10.0 |
| 621-64-7 | N-Nitroso-di-n-propylamine | 10.0 | U | 0.010 | 10.0 |
| 86-30-6 | N-Nitrosodiphenylamine | 10.0 | U | 0.010 | 10.0 |
| 95-48-7 | o-Cresol | 10.0 | U | 0.010 | 10.0 |

3/13/04
MS

SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

Lab Name: GCAL Sample ID: SKGW62A1009
 Lab Code: LA024 2 Case No.: _____ Contract: _____
 SAS No.: _____ SOG No.: 204031939 Lab File ID: S5034
 Matrix: Water Lab Sample ID: 20403220805
 Sample wt/vol: _____ Units: _____ Date Collected: 03/17/04 Time: 1150
 Level: (low/med) _____ Date Received: 03/19/04
 % Moisture: not dec. _____ Date Extracted: 3/22/04
 GC Column: RTX-SMS-30 ID: 53 (mm) Date Analyzed: 03/26/04 Time: 2112
 Concentrated Extract Volume: 1000 (µL) Dilution Factor: 1 Analyst: RLW
 Injection Volume: 1.0 (µL) Prep Method: _____
 GPC Cleanup: (Y/N) N pH: _____ Analytical Method: SW-846 8270C
 Instrument ID: MSSV2

Number TICs Found: 0

CONCENTRATION UNITS

| CAS NO. | COMPOUND | RT | EST. CONC. | Q |
|---------|------------------|----|------------|---|
| 1. | No pics detected | | | |

5/17/04
RLW

SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

Lab Name: GCAL Sample ID: SKGW641009
 Lab Code: LA024 Case No.: _____ Contract: _____
 SAS No.: _____ SDG No.: 204031909 Lab File ID: 2040326/S5035
 Matrix: Water Lab Sample ID: 20403220806
 Sample wt/vol: 1000 Units: mL Date Collected: 03/17/04 Time: 1410
 Level: (low/med) _____ Date Received: 03/19/04
 % Moisture: _____ decanted: (Y/N) _____ Date Extracted: 3/22/04
 GC Column: DB-5MS-30M ID: .25 (mm) Date Analyzed: 03/26/04 Time: 2138
 Concentrated Extract Volume: 1000 (µL) Dilution Factor: 1 Analyst: JAR3
 Injection Volume: 1.0 (µL) Prep Method: _____
 GPC Cleanup: (Y/N) N pH: _____ Analytical Method: OLMO 4.2
 Instrument ID: MSSV2

CONCENTRATION UNITS: ug/LPrep Batch: 271164 Analytical Batch: 271781

| CAS NO. | COMPOUND | RESULT | Q | MDL | PQL |
|-----------|-----------------------------|--------|---|-------|------|
| 95-95-4 | 2,4,5-Trichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 88-06-2 | 2,4,6-Trichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 120-83-2 | 2,4-Dichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 51-28-5 | 2,4-Dinitrophenol | 25.0 | U | 0.010 | 25.0 |
| 121-14-2 | 2,4-Dinitrotoluene | 10.0 | U | 0.010 | 10.0 |
| 606-20-2 | 2,6-Dinitrotoluene | 10.0 | U | 0.010 | 10.0 |
| 91-58-7 | 2-Chloronaphthalene | 10.0 | U | 0.010 | 10.0 |
| 95-57-8 | 2-Chlorophenol | 10.0 | U | 0.010 | 10.0 |
| 91-57-6 | 2-Methylnaphthalene | 10.0 | U | 0.010 | 10.0 |
| 88-74-4 | 2-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 88-75-5 | 2-Nitrophenol | 10.0 | U | 0.010 | 10.0 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 10.0 | U | 0.010 | 10.0 |
| 99-09-2 | 3-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 534-52-1 | 2-Methyl-4,6-dinitrophenol | 25.0 | U | 0.010 | 25.0 |
| 59-50-7 | 4-Chloro-3-methylphenol | 10.0 | U | 0.010 | 10.0 |
| 106-47-8 | 4-Chloroaniline | 10.0 | U | 0.010 | 10.0 |
| 7005-72-3 | 4-Chlorophenyl-phenylether | 10.0 | U | 0.010 | 10.0 |
| 106-44-5 | 4-Methylphenol (p-Cresol) | 10.0 | U | 0.010 | 10.0 |
| 83-32-9 | Acenaphthene | 10.0 | U | 0.010 | 10.0 |
| 208-96-8 | Acenaphthylene | 10.0 | U | 0.010 | 10.0 |
| 120-12-7 | Anthracene | 10.0 | U | 0.010 | 10.0 |
| 56-55-3 | Benzo(a)anthracene | 10.0 | U | 0.010 | 10.0 |
| 50-32-8 | Benzo(a)pyrene | 10.0 | U | 0.010 | 10.0 |
| 205-99-2 | Benzo(b)fluoranthene | 10.0 | U | 0.010 | 10.0 |
| 191-24-2 | Benzo(g,h,i)perylene | 10.0 | U | 0.010 | 10.0 |
| 207-08-9 | Benzo(k)fluoranthene | 10.0 | U | 0.010 | 10.0 |
| 111-91-1 | Bis(2-Chloroethoxy)methane | 10.0 | U | 0.010 | 10.0 |
| 111-44-4 | Bis(2-Chloroethyl)ether | 10.0 | U | 0.010 | 10.0 |
| 108-60-1 | bis(2-Chloroisopropyl)ether | 10.0 | U | 0.010 | 10.0 |
| 117-81-7 | bis(2-ethylhexyl)phthalate | 10.0 | U | 0.010 | 10.0 |

SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

Lab Name: GCAL Sample ID: SKGW641009
 Lab Code: LA024 Case No.: _____ Contract: _____
 SAS No.: _____ SDG No.: 204031909 Lab File ID: 2040326/S5035
 Matrix: Water Lab Sample ID: 20403220806
 Sample wt/vol: 1000 Units mL Date Collected: 03/17/04 Time: 1410
 Level: (low/med) _____ Date Received: 03/19/04
 % Moisture: _____ decanted: (Y/N) _____ Date Extracted: 3/22/04
 GC Column: DB-5MS-30M ID: 25 (mm) Date Analyzed: 03/26/04 Time: 2138
 Concentrated Extract Volume: 1000 (µL) Dilution Factor: 1 Analyst: JAR3
 Injection Volume: 1.0 (µL) Prep Method: _____
 GPC Cleanup: (Y/N) N pH: _____ Analytical Method: OLMO 4.2
 Instrument ID: MSSV2
 Prep Batch: 271164 Analytical Batch: 271781

| CAS NO. | COMPOUND | RESULT | Q | MDL | PQL |
|----------|---------------------------|--------|---|-------|------|
| 101-55-3 | 4-Bromophenyl-phenylether | 10.0 | U | 0.010 | 10.0 |
| 85-68-7 | Butylbenzylphthalate | 10.0 | U | 0.010 | 10.0 |
| 86-74-8 | Carbazole | 10.0 | U | 0.010 | 10.0 |
| 218-01-8 | Chrysene | 10.0 | U | 0.010 | 10.0 |
| 84-74-2 | Di-n-butylphthalate | 10.0 | U | 0.010 | 10.0 |
| 117-84-0 | Di-n-octylphthalate | 10.0 | U | 0.010 | 10.0 |
| 53-70-3 | Dibenz(a,h)anthracene | 10.0 | U | 0.010 | 10.0 |
| 132-64-9 | Dibenzofuran | 10.0 | U | 0.010 | 10.0 |
| 84-86-2 | Diethylphthalate | 10.0 | U | 0.010 | 10.0 |
| 131-11-3 | Dimethyl-phthalate | 10.0 | U | 0.010 | 10.0 |
| 105-67-9 | 2,4-Dimethylphenol | 10.0 | U | 0.010 | 10.0 |
| 208-44-0 | Fluoranthene | 10.0 | U | 0.010 | 10.0 |
| 86-73-7 | Fluorene | 10.0 | U | 0.010 | 10.0 |
| 118-74-1 | Hexachlorobenzene | 10.0 | U | 0.010 | 10.0 |
| 87-68-3 | Hexachlorobutadiene | 10.0 | U | 0.010 | 10.0 |
| 77-47-4 | Hexachlorocyclopentadiene | 10.0 | U | 0.010 | 10.0 |
| 87-72-1 | Hexachloroethane | 10.0 | U | 0.010 | 10.0 |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | 10.0 | U | 0.010 | 10.0 |
| 78-59-1 | Isophorone | 10.0 | U | 0.010 | 10.0 |
| 91-20-3 | Naphthalene | 10.0 | U | 0.010 | 10.0 |
| 100-01-6 | 4-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 98-95-3 | Nitrobenzene | 10.0 | U | 0.010 | 10.0 |
| 100-02-7 | 4-Nitrophenol | 25.0 | U | 0.010 | 25.0 |
| 87-86-5 | Pentachlorophenol | 25.0 | U | 0.010 | 25.0 |
| 85-01-8 | Phenanthrene | 10.0 | U | 0.010 | 10.0 |
| 108-95-2 | Phenol | 10.0 | U | 0.010 | 10.0 |
| 129-00-0 | Pyrene | 10.0 | U | 0.010 | 10.0 |
| 621-64-7 | N-Nitroso-α-n-propylamine | 10.0 | U | 0.010 | 10.0 |
| 86-30-6 | N-Nitrosodiphenylamine | 10.0 | U | 0.010 | 10.0 |
| 95-48-7 | o-Cresol | 10.0 | U | 0.010 | 10.0 |

43
45

5/13/04
mz

1F
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

| | |
|--|---|
| Lab Name: <u>GCAL</u> | Sample ID: <u>SKGW641009</u> |
| Lab Code: <u>LA024</u> <u>2</u> Case No.: _____ | Contract: _____ |
| SAS No.: _____ SDG No.: <u>204031909</u> | Lab File ID: <u>S5035</u> |
| Matrix: <u>Water</u> | Lab Sample ID: <u>20403220806</u> |
| Sample wt/vol: _____ Units: _____ | Date Collected: <u>03/17/04</u> Time: <u>1410</u> |
| Level: (low/med) _____ | Date Received: <u>03/19/04</u> |
| % Moisture: not dec. _____ | Date Extracted: <u>3/22/04</u> |
| GC Column: <u>RTX-5MS-30</u> ID: <u>.53</u> (mm) | Date Analyzed: <u>03/26/04</u> Time: <u>2206</u> |
| Concentrated Extract Volume: <u>1000</u> (µL) | Dilution Factor: <u>1</u> Analyst: <u>RLW</u> |
| Injection Volume: <u>1.0</u> (µL) | Prep Method: _____ |
| GPC Cleanup: (Y/N) <u>N</u> pH: _____ | Analytical Method: <u>SW-846 8270C</u> |
| | Instrument ID: <u>MSSV2</u> |

Number TICs Found : 0

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | RT | EST. CONC. | Q |
|---------|------------------|----|------------|---|
| 1. | No tics detected | | | |

*5/18/04
m*

SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

Lab Name: GCAL Sample ID: SKGW651009
 Lab Code: LA024 Case No. _____ Contract _____
 SAS No.: _____ SDG No.: 224031909 Lab File ID: 2040326/S5036
 Matrix: Water Lab Sample ID: 20403220807
 Sample wt/vol: 1000 Units mL Date Collected: 03/17/04 Time: 1440
 Level: (low/med) _____ Date Received: 03/19/04
 % Moisture: _____ decanted: (Y/N) _____ Date Extracted: 3/22/04
 GC Column: DB-5MS-30M ID: 25 (mm) Date Analyzed: 03/26/04 Time: 2206
 Concentrated Extract Volume: 1000 (µL) Dilution Factor: 1 Analyst: JAR3
 Injection Volume: 1.0 (µL) Prep Method: _____
 GPC Cleanup: (Y/N) N pH: _____ Analytical Method: OLMO 4.2
 Instrument ID: MSSV2
 Prep Batch: 271164 Analytical Batch: 271781

| CAS NO. | COMPOUND | RESULT | Q | MDL | PQL |
|-----------|-----------------------------|--------|---|-------|------|
| 95-95-4 | 2,4,5-Trichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 88-06-2 | 2,4,6-Trichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 120-83-2 | 2,4-Dichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 51-28-5 | 2,4-Dinitrophenol | 25.0 | U | 0.010 | 25.0 |
| 121-14-2 | 2,4-Dinitrotoluene | 10.0 | U | 0.010 | 10.0 |
| 606-20-2 | 2,6-Dinitrotoluene | 10.0 | U | 0.010 | 10.0 |
| 91-58-7 | 2-Chloronaphthalene | 10.0 | U | 0.010 | 10.0 |
| 95-57-8 | 2-Chlorophenol | 10.0 | U | 0.010 | 10.0 |
| 91-57-6 | 2-Methylnaphthalene | 10.0 | U | 0.010 | 10.0 |
| 88-74-4 | 2-Nitroanisole | 25.0 | U | 0.010 | 25.0 |
| 88-75-5 | 2-Nitrophenol | 10.0 | U | 0.010 | 10.0 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 10.0 | U | 0.010 | 10.0 |
| 99-09-2 | 3-Nitroanisole | 25.0 | U | 0.010 | 25.0 |
| 534-52-1 | 2-Methyl-4,6-dinitrophenol | 25.0 | U | 0.010 | 25.0 |
| 59-50-7 | 4-Chloro-3-methylphenol | 10.0 | U | 0.010 | 10.0 |
| 106-47-8 | 4-Chloroaniline | 10.0 | U | 0.010 | 10.0 |
| 7005-72-3 | 4-Chlorophenyl-phenylether | 10.0 | U | 0.010 | 10.0 |
| 106-44-5 | 4-Methylphenol (p-Cresol) | 10.0 | U | 0.010 | 10.0 |
| 83-32-9 | Acenaphthene | 10.0 | U | 0.010 | 10.0 |
| 208-96-8 | Acenaphthylene | 10.0 | U | 0.010 | 10.0 |
| 120-12-7 | Anthracene | 10.0 | U | 0.010 | 10.0 |
| 56-55-3 | Benzo(a)anthracene | 10.0 | U | 0.010 | 10.0 |
| 50-32-6 | Benzo(a)pyrene | 10.0 | U | 0.010 | 10.0 |
| 205-99-2 | Benzo(b)fluoranthene | 10.0 | U | 0.010 | 10.0 |
| 191-24-2 | Benzo(g,h,i)perylene | 10.0 | U | 0.010 | 10.0 |
| 207-08-9 | Benzo(k)fluoranthene | 10.0 | U | 0.010 | 10.0 |
| 111-91-1 | Bis(2-Chloroethoxy)methane | 10.0 | U | 0.010 | 10.0 |
| 111-44-4 | Bis(2-Chloroethyl)ether | 10.0 | U | 0.010 | 10.0 |
| 108-60-1 | bis(2-Chloroisopropyl)ether | 10.0 | U | 0.010 | 10.0 |
| 117-81-7 | bis(2-ethylhexyl)phthalate | 10.0 | U | 0.010 | 10.0 |

SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

Lab Name: GCALSample ID: SKGW651009Lab Code: LA024 Case No.: _____

Contract: _____

SAS No.: _____ SDG No.: 204031909Lab File ID: 2040326/S5036Matrix: WaterLab Sample ID: 20403220807Sample wt/vol: 1000 Units: mLDate Collected: 03/17/04 Time: 1440

Level: (low/med) _____

Date Received: 03/19/04

% Moisture: _____ decanted: (Y/N) _____

Date Extracted: 3/21/04GC Column: DB-5MS-30M ID: .25 (mm)Date Analyzed: 03/26/04 Time: 2206Concentrated Extract Volume: 1000 (μL)Dilution Factor: 1 Analyst: JAR3Injection Volume: 1.0 (μL)

Prep Method: _____

GPC Cleanup: (Y/N) N pH: _____Analytical Method: OLMO 4.2Instrument ID: MSSV2CONCENTRATION UNITS: ug/LPrep Batch: 271164 Analytical Batch: 271781

| CAS NO. | COMPOUND | RESULT | Q | MDL | PQL |
|----------|----------------------------|--------|---|-------|------|
| 101-55-3 | 4-Bromophenyl-phenylether | 10.0 | U | 0.010 | 10.0 |
| 85-68-7 | Butylbenzylphthalate | 10.0 | U | 0.010 | 10.0 |
| 86-74-8 | Carbazole | 10.0 | U | 0.010 | 10.0 |
| 218-01-9 | Chrysene | 10.0 | U | 0.010 | 10.0 |
| 84-74-2 | Di-n-butylphthalate | 10.0 | U | 0.010 | 10.0 |
| 117-84-0 | Di-n-octylphthalate | 10.0 | U | 0.010 | 10.0 |
| 53-70-3 | Dibenz(a,h)anthracene | 10.0 | U | 0.010 | 10.0 |
| 132-64-9 | Dibenzofuran | 10.0 | U | 0.010 | 10.0 |
| 84-66-2 | Diethylphthalate | 10.0 | U | 0.010 | 10.0 |
| 131-11-3 | Dimethyl-phthalate | 10.0 | U | 0.010 | 10.0 |
| 105-67-9 | 2,4-Dimethylphenol | 10.0 | U | 0.010 | 10.0 |
| 206-44-0 | Fluoranthene | 10.0 | U | 0.010 | 10.0 |
| 86-73-7 | Fluorene | 10.0 | U | 0.010 | 10.0 |
| 118-74-1 | Hexachlorobenzene | 10.0 | U | 0.010 | 10.0 |
| 87-68-3 | Hexachlorobutadiene | 10.0 | U | 0.010 | 10.0 |
| 77-47-4 | Hexachlorocyclopentadiene | 10.0 | U | 0.010 | 10.0 |
| 67-72-1 | Hexachloroethane | 10.0 | U | 0.010 | 10.0 |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | 10.0 | U | 0.010 | 10.0 |
| 78-59-1 | Isophorone | 10.0 | U | 0.010 | 10.0 |
| 91-20-3 | Naphthalene | 10.0 | U | 0.010 | 10.0 |
| 100-01-6 | 4-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 98-95-3 | Nitrobenzene | 10.0 | U | 0.010 | 10.0 |
| 100-02-7 | 4-Nitrophenol | 25.0 | U | 0.010 | 25.0 |
| 87-86-5 | Pentachlorophenol | 25.0 | U | 0.010 | 25.0 |
| 85-01-8 | Phenanthrene | 10.0 | U | 0.010 | 10.0 |
| 108-95-2 | Phenol | 10.0 | U | 0.010 | 10.0 |
| 129-00-0 | Pyrene | 10.0 | U | 0.010 | 10.0 |
| 621-64-7 | N-Nitroso-di-n-propylamine | 10.0 | U | 0.010 | 10.0 |
| 86-30-6 | N-Nitrosodiphenylamine | 10.0 | U | 0.010 | 10.0 |
| 95-48-7 | o-Cresol | 10.0 | U | 0.010 | 10.0 |

WS
WS5/13/04
MSE

SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

| | |
|--|---|
| Lab Name: <u>GCAL</u> | Sample ID: <u>SKGW651009</u> |
| Lab Code: <u>LA024</u> <u>2</u> Case No.: _____ | Contract: _____ |
| SAS No.: _____ SDG No.: <u>204031909</u> | Lab File ID: <u>S5036</u> |
| Matrix: <u>Water</u> | Lab Sample ID: <u>20403220807</u> |
| Sample wt/vol: _____ Units: _____ | Date Collected: <u>03/17/04</u> Time: <u>1440</u> |
| Level: (low/med) _____ | Date Received: <u>03/19/04</u> |
| % Moisture: not dec. _____ | Date Extracted: <u>3/22/04</u> |
| GC Column: <u>RTX-5MS-30</u> ID: <u>.53</u> (mm) | Date Analyzed: <u>03/26/04</u> Time: <u>2206</u> |
| Concentrated Extract Volume: <u>1000</u> (µL) | Dilution Factor: <u>1</u> Analyst: <u>RLW</u> |
| Injection Volume: <u>1.0</u> (µL) | Prep Method: _____ |
| GPC Cleanup: (Y/N) <u>N</u> pH: _____ | Analytical Method: <u>SW-846 8270C</u> |
| | Instrument ID: <u>MSSV2</u> |

Number TICs Found: 0

CONCENTRATION UNITS

| CAS NO. | COMPOUND | RT | EST. CONC. | Q |
|---------|------------------|----|------------|---|
| 1. | No tics detected | | | |

5/17/04
mx

1D
ORGANICS ANALYSIS DATA SHEET

Lab Name: GCAL Sample ID: SKGW581009
 Lab Code: LA024 Case No.: _____ Contract: _____
 Matrix: Water SAS No.: _____ SDG No.: 204031909
 Sample wt/vol: 1000 Units: mL Lab Sample ID: 20403190901
 Level: (low/med) _____ Date Collected: 03/18/04 Time: 1148
 % Moisture: _____ decanted: (Y/N) _____ Date Received: 03/19/04
 GC Column: DB-608-30M ID: .53 (mm) Date Extracted: 3/23/04
 Concentrated Extract Volume: 1000 (µL) Date Analyzed: 04/01/04 Time: 0502
 Soil Aliquot Volume: _____ (µL) Dilution Factor: 1 Analyst: TLS
 Injection Volume: 1 (µL) Prep Method: _____
 GPC Cleanup: (Y/N) N pH: _____ Analytical Method: OLMO 4.2
 Prep Batch: 271129 Analytical Batch: 271697 Sulfur Cleanup: (Y/N) N Instrument ID: GCS6A
 Lab File ID: 2040331/SV6024

CONCENTRATION UNITS: ug/L

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|------------|---------------------|--------|---|---------|-------|
| 72-54-8 | 4,4'-DDD | 0.100 | U | 0.00010 | 0.100 |
| 72-55-9 | 4,4'-DDE | 0.100 | U | 0.00010 | 0.100 |
| 50-29-3 | 4,4'-DDT | 0.100 | U | 0.00010 | 0.100 |
| 309-00-2 | Aldrin | 0.050 | U | 0.00010 | 0.050 |
| 12674-11-2 | Aroclor-1016 | 1.00 | U | 0.00010 | 1.00 |
| 11104-28-2 | Aroclor-1221 | 2.00 | U | 0.00010 | 2.00 |
| 11141-16-5 | Aroclor-1232 | 1.00 | U | 0.00010 | 1.00 |
| 53469-21-9 | Aroclor-1242 | 1.00 | U | 0.00010 | 1.00 |
| 12672-29-6 | Aroclor-1248 | 1.00 | U | 0.00010 | 1.00 |
| 11097-69-1 | Aroclor-1254 | 1.00 | U | 0.00010 | 1.00 |
| 11096-82-5 | Aroclor-1260 | 1.00 | U | 0.00010 | 1.00 |
| 60-57-1 | Dieldrin | 0.100 | U | 0.00010 | 0.100 |
| 959-98-8 | Endosulfan I | 0.050 | U | 0.00010 | 0.050 |
| 33213-85-9 | Endosulfan II | 0.100 | U | 0.00010 | 0.100 |
| 1031-07-8 | Endosulfan sulfate | 0.100 | U | 0.00010 | 0.100 |
| 72-20-8 | Endrin | 0.100 | U | 0.00010 | 0.100 |
| 7421-93-4 | Endrin aldehyde | 0.100 | U | 0.00010 | 0.100 |
| 53494-70-5 | Endrin ketone | 0.100 | U | 0.00010 | 0.100 |
| 76-44-8 | Heptachlor | 0.050 | U | 0.00010 | 0.050 |
| 1024-57-3 | Heptachlor epoxide | 0.050 | U | 0.00010 | 0.050 |
| 72-43-5 | Methoxychlor | 0.500 | U | 0.00010 | 0.500 |
| 8001-35-2 | Toxaphene | 5.00 | U | 0.00010 | 5.00 |
| 319-84-6 | alpha-BHC | 0.050 | U | 0.00010 | 0.050 |
| 5103-71-9 | alpha-Chlordane | 0.050 | U | 0.00010 | 0.050 |
| 319-85-7 | beta-BHC | 0.050 | U | 0.00010 | 0.050 |
| 319-86-8 | delta-BHC | 0.050 | U | 0.00010 | 0.050 |
| 58-89-9 | gamma-BHC (Lindane) | 0.050 | U | 0.00010 | 0.050 |
| 5103-74-2 | gamma-Chlordane | 0.050 | U | 0.00010 | 0.050 |

ORGANICS ANALYSIS DATA SHEET

Lab Name: GCAL Sample ID: SKGW58DUP1009
 Lab Code: LA024 Case No.: _____ Contract: _____
 Matrix: Water SAS No: _____ SDG No.: 204031909
 Sample wt/vol: 1000 Units mL Lab Sample ID: 20403190902
 Level: (low/med) _____ Date Collected: 03/18/04 Time: 1208
 % Moisture: _____ decanted: (Y/N) _____ Date Received: 03/19/04
 GC Column: DB-608-30M ID: 53 (mm) Date Extracted: 3/23/04
 Concentrated Extract Volume: 1000 (µL) Date Analyzed: 04/01/04 Time: 0530
 Soil Aliquot Volume: _____ (µL) Dilutor Factor: 1 Analyst: TLS
 Injection Volume: 1 (µL) Prep Method: _____
 GPC Cleanup: (Y/N) N pH: _____ Analytical Method: OLMO 4.2
 Prep Batch: 271129 Analytical Batch: 271697 Sulfur Cleanup: (Y/N) N Instrument ID: GCS6A
 Lab File ID: 2040331/SV6025

CONCENTRATION UNITS: µg/L

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|------------|---------------------|--------|---|---------|-------|
| 72-54-8 | 4,4'-DDD | 0.100 | U | 0.00010 | 0.100 |
| 72-55-9 | 4,4'-DDE | 0.100 | U | 0.00010 | 0.100 |
| 50-29-3 | 4,4'-DDT | 0.100 | U | 0.00010 | 0.100 |
| 309-00-2 | Aldrin | 0.050 | U | 0.00010 | 0.050 |
| 12874-11-2 | Aroclor-1016 | 1.00 | U | 0.00010 | 1.00 |
| 11104-28-2 | Aroclor-1221 | 2.00 | U | 0.00010 | 2.00 |
| 11141-16-5 | Aroclor-1232 | 1.00 | U | 0.00010 | 1.00 |
| 53489-21-9 | Aroclor-1242 | 1.00 | U | 0.00010 | 1.00 |
| 12872-29-6 | Aroclor-1248 | 1.00 | U | 0.00010 | 1.00 |
| 11097-69-1 | Aroclor-1254 | 1.00 | U | 0.00010 | 1.00 |
| 11096-82-5 | Aroclor-1260 | 1.00 | U | 0.00010 | 1.00 |
| 60-57-1 | Dieldrin | 0.100 | U | 0.00010 | 0.100 |
| 959-98-8 | Endosulfan I | 0.050 | U | 0.00010 | 0.050 |
| 33213-65-9 | Endosulfan II | 0.100 | U | 0.00010 | 0.100 |
| 1031-07-8 | Endosulfan sulfate | 0.100 | U | 0.00010 | 0.100 |
| 72-20-8 | Endrin | 0.100 | U | 0.00010 | 0.100 |
| 7421-93-4 | Endrin aldehyde | 0.100 | U | 0.00010 | 0.100 |
| 53494-70-5 | Endrin ketone | 0.100 | U | 0.00010 | 0.100 |
| 76-44-8 | Heptachlor | 0.050 | U | 0.00010 | 0.050 |
| 1024-57-3 | Heptachlor epoxide | 0.050 | U | 0.00010 | 0.050 |
| 72-43-5 | Methoxychlor | 0.500 | U | 0.00010 | 0.500 |
| 8001-35-2 | Toxaphene | 5.00 | U | 0.00010 | 5.00 |
| 319-84-6 | alpha-BHC | 0.050 | U | 0.00010 | 0.050 |
| 5103-71-9 | alpha-Chlordane | 0.050 | U | 0.00010 | 0.050 |
| 319-85-7 | beta-BHC | 0.050 | U | 0.00010 | 0.050 |
| 319-86-8 | delta-BHC | 0.050 | U | 0.00010 | 0.050 |
| 58-89-9 | gamma-BHC (Lindane) | 0.050 | U | 0.00010 | 0.050 |
| 5103-74-2 | gamma-Chlordane | 0.050 | U | 0.00010 | 0.050 |

1D
ORGANICS ANALYSIS DATA SHEET

| | |
|---|--|
| Lab Name: <u>GCAL</u> | Sample ID: <u>SKGW631009</u> |
| Lab Code: <u>LA024</u> Case No.: _____ | Contract: _____ |
| Matrix: <u>Water</u> | SAS No.: _____ SDG No.: <u>204031909</u> |
| Sample wt/vol: <u>1000</u> Units: <u>mL</u> | Lab Sample ID: <u>20403190903</u> |
| Level: (low/med) _____ | Date Collected: <u>03/18/04</u> Time: <u>1430</u> |
| % Moisture: _____ decanted: (Y/N) _____ | Date Received: <u>03/19/04</u> |
| GC Column: <u>DB-608-30M</u> ID: <u>.53</u> (mm) | Date Extracted: <u>3/23/04</u> |
| Concentrated Extract Volume: <u>1000</u> (µL) | Date Analyzed: <u>04/01/04</u> Time: <u>0558</u> |
| Soil Aliquot Volume: _____ (µL) | Dilution Factor: <u>1</u> Analyst: <u>TLS</u> |
| Injection Volume: <u>1</u> (µL) | Prep Method: _____ |
| GPC Cleanup: (Y/N) <u>N</u> pH: _____ | Analytical Method: <u>OLMO 4.2</u> |
| Prep Batch: <u>271129</u> Analytical Batch: <u>271697</u> | Sulfur Cleanup: (Y/N) <u>N</u> Instrument ID: <u>GCS6A</u> |
| CONCENTRATION UNITS: <u>ug/L</u> | Lab File ID: <u>2040331/SV6026</u> |

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|------------|---------------------|--------|---|---------|-------|
| 72-54-8 | 4,4'-DDD | 0.100 | U | 0.00010 | 0.100 |
| 72-55-9 | 4,4'-DDE | 0.100 | U | 0.00010 | 0.100 |
| 50-29-3 | 4,4'-DDT | 0.100 | U | 0.00010 | 0.100 |
| 309-00-2 | Aldrin | 0.050 | U | 0.00010 | 0.050 |
| 12674-11-2 | Aroclor-1016 | 1.00 | U | 0.00010 | 1.00 |
| 11104-28-2 | Aroclor-1221 | 2.00 | U | 0.00010 | 2.00 |
| 11141-16-5 | Aroclor-1232 | 1.00 | U | 0.00010 | 1.00 |
| 53469-21-9 | Aroclor-1242 | 1.00 | U | 0.00010 | 1.00 |
| 12672-29-6 | Aroclor-1248 | 1.00 | U | 0.00010 | 1.00 |
| 11097-69-1 | Aroclor-1254 | 1.00 | U | 0.00010 | 1.00 |
| 11096-82-5 | Aroclor-1260 | 1.00 | U | 0.00010 | 1.00 |
| 60-57-1 | Dieldrin | 0.100 | U | 0.00010 | 0.100 |
| 959-98-8 | Endosulfan I | 0.050 | U | 0.00010 | 0.050 |
| 33213-65-9 | Endosulfan II | 0.100 | U | 0.00010 | 0.100 |
| 1031-07-8 | Endosulfan sulfate | 0.100 | U | 0.00010 | 0.100 |
| 72-20-8 | Endrin | 0.100 | U | 0.00010 | 0.100 |
| 7421-93-4 | Endrin aldehyde | 0.100 | U | 0.00010 | 0.100 |
| 53494-70-6 | Endrin ketone | 0.100 | U | 0.00010 | 0.100 |
| 78-44-8 | Endrin epoxide | 0.050 | U | 0.00010 | 0.050 |
| 1024-47-3 | Endrin epoxide | 0.050 | U | 0.00010 | 0.050 |
| 72-43-5 | Endrin epoxide | 0.500 | U | 0.00010 | 0.500 |
| 8001-35-2 | Toxaphene | 5.00 | U | 0.00010 | 5.00 |
| 319-84-6 | alpha-BHC | 0.050 | U | 0.00010 | 0.050 |
| 5103-71-9 | alpha-Chlordane | 0.050 | U | 0.00010 | 0.050 |
| 319-85-7 | beta-BHC | 0.050 | U | 0.00010 | 0.050 |
| 319-86-8 | delta-BHC | 0.050 | U | 0.00010 | 0.050 |
| 58-89-9 | gamma-BHC (Lindane) | 0.050 | U | 0.00010 | 0.050 |
| 5103-74-2 | gamma-Chlordane | 0.050 | U | 0.00010 | 0.050 |

ORGANICS ANALYSIS DATA SHEET

Lab Name: GCAL Sample ID: SKGW611009

Lab Code: LA024 Case No.: _____ Contract: _____

Matrix: Water SAS No: _____ SDG No.: 204031909

Sample wt/vol: 1000 Units: mL Lab Sample ID: 20403190904

Level: (low/mid) _____ Date Collected: 03/18/04 Time: 1535

% Moisture: _____ decanted: (Y/N) _____ Date Received: 03/19/04

GC Column: DB-608-30M ID: .53 (mm) Date Extracted: 3/23/04

Concentrated Extract Volume: 1000 (µL) Date Analyzed: 04/01/04 Time: 0627

Soil Aliquot Volume: _____ (µL) Dilution Factor: 1 Analyst: TLS

Injection Volume: 1 (µL) Prep Method: _____

GPC Cleanup: (Y/N) N pH: _____ Analytical Method: OLMO 4.2

Prep Batch: 271129 Analytical Batch: 271697 Sulfur Cleanup: (Y/N) N Instrument ID: GCS6A

Lab File ID: 2040331/SV6027

CONCENTRATION UNITS: µg/L

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|------------|---------------------|--------|---|---------|-------|
| 72-54-8 | 4,4'-DDD | 0.100 | U | 0.00010 | 0.100 |
| 72-55-9 | 4,4'-DDE | 0.100 | U | 0.00010 | 0.100 |
| 50-29-3 | 4,4'-DDT | 0.100 | U | 0.00010 | 0.100 |
| 309-00-2 | Aldrin | 0.050 | U | 0.00010 | 0.050 |
| 12674-11-2 | Aroclor-1016 | 1.00 | U | 0.00010 | 1.00 |
| 11104-28-2 | Aroclor-1221 | 2.00 | U | 0.00010 | 2.00 |
| 11141-16-5 | Aroclor-1232 | 1.00 | U | 0.00010 | 1.00 |
| 53469-21-9 | Aroclor-1242 | 1.00 | U | 0.00010 | 1.00 |
| 12672-29-6 | Aroclor-1248 | 1.00 | U | 0.00010 | 1.00 |
| 11097-89-1 | Aroclor-1254 | 1.00 | U | 0.00010 | 1.00 |
| 11096-82-5 | Aroclor-1260 | 1.00 | U | 0.00010 | 1.00 |
| 60-57-1 | Octadrin | 0.100 | U | 0.00010 | 0.100 |
| 959-98-8 | Endosulfan I | 0.050 | U | 0.00010 | 0.050 |
| 33213-65-9 | Endosulfan II | 0.100 | U | 0.00010 | 0.100 |
| 1031-07-8 | Endosulfan sulfate | 0.100 | U | 0.00010 | 0.100 |
| 72-20-8 | Endrin | 0.100 | U | 0.00010 | 0.100 |
| 7421-93-4 | Endrin aldehyde | 0.100 | U | 0.00010 | 0.100 |
| 53494-70-5 | Endrin ketone | 0.100 | U | 0.00010 | 0.100 |
| 76-44-8 | Heptachlor | 0.050 | U | 0.00010 | 0.050 |
| 1024-57-3 | Heptachlor epoxide | 0.050 | U | 0.00010 | 0.050 |
| 72-43-5 | Methoxychlor | 0.500 | U | 0.00010 | 0.500 |
| 8001-35-2 | Toxaphene | 5.00 | U | 0.00010 | 5.00 |
| 319-84-6 | alpha-BHC | 0.050 | U | 0.00010 | 0.050 |
| 5103-71-9 | alpha-Chlordane | 0.050 | U | 0.00010 | 0.050 |
| 319-85-7 | beta-BHC | 0.050 | U | 0.00010 | 0.050 |
| 319-86-8 | delta-BHC | 0.050 | U | 0.00010 | 0.050 |
| 58-89-9 | gamma-BHC (lincane) | 0.050 | U | 0.00010 | 0.050 |
| 5103-74-2 | gamma-Chlordane | 0.050 | U | 0.00010 | 0.050 |

5/24/04
msc

1D
ORGANICS ANALYSIS DATA SHEET

Lab Name: GCAL Sample ID: SKGWFB1009
 Lab Code: LA024 Case No.: _____ Contract: _____
 Matrix: Water SAS No.: _____ SDG No.: 204031909
 Sample wt/vol: 1000 Units: mL Lab Sample ID: 20403190908
 Level: (low/med) _____ Date Collected: 03/18/04 Time: 1702
 % Moisture: _____ decanted: (Y/N) _____ Date Received: 03/19/04
 GC Column: DB-608-30M ID: .53 (mm) Date Extracted: 3/23/04
 Concentrated Extract Volume: 1000 (µL) Date Analyzed: 04/01/04 Time: 0916
 Soil Aliquot Volume: _____ (µL) Dilution Factor: 1 Analyst: TLS
 Injection Volume: 1 (µL) Prep Method: _____
 GPC Cleanup: (Y/N) N pH: _____ Analytical Method: OLMO 4.2
 Prep Batch: 271129 Analytical Batch: 271697 Sulfur Cleanup: (Y/N) N Instrument ID: GCS6A
 CONCENTRATION UNITS: ug/L Lab File ID: 2040331/SV6033

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|------------|---------------------|--------|---|---------|-------|
| 72-54-8 | 4,4'-DDD | 0.100 | U | 0.00010 | 0.100 |
| 72-55-9 | 4,4'-DDE | 0.100 | U | 0.00010 | 0.100 |
| 50-29-3 | 4,4'-DDT | 0.100 | U | 0.00010 | 0.100 |
| 309-00-2 | Aldrin | 0.050 | U | 0.00010 | 0.050 |
| 12674-11-2 | Aroclor-1016 | 1.00 | U | 0.00010 | 1.00 |
| 11104-28-2 | Aroclor-1221 | 2.00 | U | 0.00010 | 2.00 |
| 11141-16-5 | Aroclor-1232 | 1.00 | U | 0.00010 | 1.00 |
| 53469-21-9 | Aroclor-1242 | 1.00 | U | 0.00010 | 1.00 |
| 12672-29-6 | Aroclor-1248 | 1.00 | U | 0.00010 | 1.00 |
| 11097-69-1 | Aroclor-1254 | 1.00 | U | 0.00010 | 1.00 |
| 11096-82-5 | Aroclor-1260 | 1.00 | U | 0.00010 | 1.00 |
| 60-57-1 | Dieldrin | 0.100 | U | 0.00010 | 0.100 |
| 959-98-8 | Endosulfan I | 0.050 | U | 0.00010 | 0.050 |
| 33213-65-9 | Endosulfan II | 0.100 | U | 0.00010 | 0.100 |
| 1031-07-8 | Endosulfan sulfate | 0.100 | U | 0.00010 | 0.100 |
| 72-20-8 | Endrin | 0.100 | U | 0.00010 | 0.100 |
| 7421-93-4 | Endrin aldehyde | 0.100 | U | 0.00010 | 0.100 |
| 53494-70-5 | Endrin ketone | 0.100 | U | 0.00010 | 0.100 |
| 76-44-8 | Heptachlor | 0.050 | U | 0.00010 | 0.050 |
| 1024-57-3 | Heptachlor epoxide | 0.050 | U | 0.00010 | 0.050 |
| 72-43-5 | Methoxychlor | 0.500 | U | 0.00010 | 0.500 |
| 8001-35-2 | Toxaphene | 5.00 | U | 0.00010 | 5.00 |
| 319-84-6 | alpha-BHC | 0.050 | U | 0.00010 | 0.050 |
| 5103-71-9 | alpha-Chlordane | 0.050 | U | 0.00010 | 0.050 |
| 319-85-7 | beta-BHC | 0.050 | U | 0.00010 | 0.050 |
| 319-86-8 | delta-BHC | 0.050 | U | 0.00010 | 0.050 |
| 58-89-9 | gamma-BHC (Lindane) | 0.050 | U | 0.00010 | 0.050 |
| 5103-74-2 | gamma-Chlordane | 0.050 | U | 0.00010 | 0.050 |

ORGANICS ANALYSIS DATA SHEET

| | |
|---|--|
| Lab Name: <u>GCAL</u> | Sample ID: <u>SKGW06R1009</u> |
| Lab Code: <u>LA024</u> Case No.: _____ | Contract: _____ |
| Matrix: <u>Water</u> | SAS No: _____ SDG No.: <u>204031909</u> |
| Sample wt/vol: <u>1000</u> Units: <u>mL</u> | Lab Sample ID: <u>20403220801</u> |
| Level: (low/med) _____ | Date Collected: <u>03/16/04</u> Time: <u>1458</u> |
| % Moisture: _____ decanted: (Y/N) _____ | Date Received: <u>03/19/04</u> |
| GC Column: <u>DB-608-30M</u> ID: <u>.53</u> (mm) | Date Extracted: <u>3/23/04</u> |
| Concentrated Extract Volume: <u>1000</u> (µL) | Date Analyzed: <u>04/01/04</u> Time: <u>0405</u> |
| Soil Aliquot Volume: _____ (µL) | Dilution Factor: <u>1</u> Analyst: <u>TLS</u> |
| Injection Volume: <u>1</u> (µL) | Prep Method: _____ |
| GPC Cleanup: (Y/N) <u>N</u> pH: _____ | Analytical Method: <u>OLMO 4.2</u> |
| Prep Batch: <u>271129</u> Analytical Batch: <u>271697</u> | Sulfur Cleanup: (Y/N) <u>N</u> Instrument ID: <u>GCS6A</u> |
| CONCENTRATION UNITS: <u>ug/L</u> | Lab File ID: <u>2040331/SV6022</u> |

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|------------|---------------------|--------|---|---------|-------|
| 72-54-8 | 4,4'-DDD | 0.100 | U | 0.00010 | 0.100 |
| 72-55-9 | 4,4'-DDE | 0.100 | U | 0.00010 | 0.100 |
| 50-29-3 | 4,4'-DDT | 0.100 | U | 0.00010 | 0.100 |
| 309-00-2 | Aldrin | 0.050 | U | 0.00010 | 0.050 |
| 12674-11-2 | Aroclor-1018 | 1.00 | U | 0.00010 | 1.00 |
| 11104-28-2 | Aroclor-1221 | 2.00 | U | 0.00010 | 2.00 |
| 11141-16-5 | Aroclor-1232 | 1.00 | U | 0.00010 | 1.00 |
| 53489-21-8 | Aroclor-1242 | 1.00 | U | 0.00010 | 1.00 |
| 12672-29-6 | Aroclor-1248 | 1.00 | U | 0.00010 | 1.00 |
| 11087-69-1 | Aroclor-1254 | 1.00 | U | 0.00010 | 1.00 |
| 11096-82-5 | Aroclor-1260 | 1.00 | U | 0.00010 | 1.00 |
| 80-57-1 | Dieldrin | 0.100 | U | 0.00010 | 0.100 |
| 959-98-8 | Endosulfan I | 0.050 | U | 0.00010 | 0.050 |
| 33213-65-9 | Endosulfan II | 0.100 | U | 0.00010 | 0.100 |
| 1031-07-8 | Endosulfan sulfate | 0.100 | U | 0.00010 | 0.100 |
| 72-20-8 | Endrin | 0.100 | U | 0.00010 | 0.100 |
| 7421-93-4 | Endrin aldehyde | 0.100 | U | 0.00010 | 0.100 |
| 53484-70-5 | Endrin ketone | 0.100 | U | 0.00010 | 0.100 |
| 78-44-8 | Heptachlor | 0.050 | U | 0.00010 | 0.050 |
| 1024-57-3 | Heptachlor epoxide | 0.050 | U | 0.00010 | 0.050 |
| 72-43-5 | Methoxychlor | 0.500 | U | 0.00010 | 0.500 |
| 8001-35-2 | Toxaphene | 5.00 | U | 0.00010 | 5.00 |
| 319-84-6 | alpha-BHC | 0.050 | U | 0.00010 | 0.050 |
| 5103-71-9 | alpha-Chlordane | 0.050 | U | 0.00010 | 0.050 |
| 319-85-7 | beta-BHC | 0.050 | U | 0.00010 | 0.050 |
| 319-86-8 | delta-BHC | 0.050 | U | 0.00010 | 0.050 |
| 58-89-9 | gamma-BHC (Lincane) | 0.050 | U | 0.00010 | 0.050 |
| 5103-74-2 | gamma-Chlordane | 0.050 | U | 0.00010 | 0.050 |

1D
ORGANICS ANALYSIS DATA SHEET

Lab Name: GCAL Sample ID: SKGW07R1009
 Lab Code: LA024 Case No.: _____ Contract: _____
 Matrix: Water SAS No.: _____ SDG No.: 204031909
 Sample wt/vol: 1000 Units: mL Lab Sample ID: 20403220802
 Level: (low/med) _____ Date Collected: 03/16/04 Time: 1552
 % Moisture: _____ decanted: (Y/N) _____ Date Received: 03/19/04
 GC Column: DB-608-30M ID: .53 (mm) Date Extracted: 3/23/04
 Concentrated Extract Volume: 1000 (µL) Date Analyzed: 04/01/04 Time: 0434
 Soil Aliquot Volume: _____ (µL) Dilution Factor: 1 Analyst: TLS
 Injection Volume: 1 (µL) Prep Method: _____
 GPC Cleanup: (Y/N) N pH: _____ Analytical Method: OLMO 4.2
 Prep Batch: 271129 Analytical Batch: 271697 Sulfur Cleanup: (Y/N) N Instrument ID: GCS6A
 Lab File ID: 2040331/SV6023

CONCENTRATION UNITS: ug/L

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|------------|---------------------|--------|---|---------|-------|
| 72-54-8 | 4,4'-DDD | 0.100 | U | 0.00010 | 0.100 |
| 72-55-9 | 4,4'-DDE | 0.100 | U | 0.00010 | 0.100 |
| 50-29-3 | 4,4'-DDT | 0.100 | U | 0.00010 | 0.100 |
| 309-00-2 | Aldrin | 0.050 | U | 0.00010 | 0.050 |
| 12674-11-2 | Aroclor-1016 | 1.00 | U | 0.00010 | 1.00 |
| 11104-28-2 | Aroclor-1221 | 2.00 | U | 0.00010 | 2.00 |
| 11141-16-5 | Aroclor-1232 | 1.00 | U | 0.00010 | 1.00 |
| 53469-21-9 | Aroclor-1242 | 1.00 | U | 0.00010 | 1.00 |
| 12672-29-6 | Aroclor-1248 | 1.00 | U | 0.00010 | 1.00 |
| 11097-69-1 | Aroclor-1254 | 1.00 | U | 0.00010 | 1.00 |
| 11096-82-5 | Aroclor-1260 | 1.00 | U | 0.00010 | 1.00 |
| 60-57-1 | Dieldrin | 0.100 | U | 0.00010 | 0.100 |
| 959-98-8 | Endosulfan I | 0.050 | U | 0.00010 | 0.050 |
| 33213-65-9 | Endosulfan II | 0.100 | U | 0.00010 | 0.100 |
| 1031-07-8 | Endosulfan sulfate | 0.100 | U | 0.00010 | 0.100 |
| 72-20-8 | Endrin | 0.100 | U | 0.00010 | 0.100 |
| 7421-93-4 | Endrin aldehyde | 0.100 | U | 0.00010 | 0.100 |
| 53494-70-5 | Endrin ketone | 0.100 | U | 0.00010 | 0.100 |
| 76-44-8 | Heptachlor | 0.050 | U | 0.00010 | 0.050 |
| 1024-57-3 | Heptachlor epoxide | 0.050 | U | 0.00010 | 0.050 |
| 72-43-5 | Methoxychlor | 0.500 | U | 0.00010 | 0.500 |
| 8001-35-2 | Toxaphene | 5.00 | U | 0.00010 | 5.00 |
| 319-84-6 | alpha-BHC | 0.050 | U | 0.00010 | 0.050 |
| 5103-71-9 | alpha-Chlordane | 0.050 | U | 0.00010 | 0.050 |
| 319-85-7 | beta-BHC | 0.050 | U | 0.00010 | 0.050 |
| 319-86-8 | delta-BHC | 0.050 | U | 0.00010 | 0.050 |
| 58-89-9 | gamma-BHC (Lindane) | 0.050 | U | 0.00010 | 0.050 |
| 5103-74-2 | gamma-Chlordane | 0.050 | U | 0.00010 | 0.050 |

ORGANICS ANALYSIS DATA SHEET

Lab Name: GCAL Sample ID: SKGW591009
 Lab Code: LA024 Case No.: _____ Contract: _____
 Matrix: Water SAS No. _____ SDG No.: 204031909
 Sample wt/vol: 1000 Units: mL Lab Sample ID: 20403220003
 Level: (low/mid) _____ Date Collected: 03/17/04 Time: 1050
 % Moisture: _____ decanted: (Y/N) _____ Date Received: 03/19/04
 GC Column: DB-608-30M ID: 53 (mm) Date Extracted: 3/23/04
 Concentrated Extract Volume: 1000 (µL) Date Analyzed: 04/01/04 Time: 0944
 Soil Aliquot Volume: _____ (µL) Dilution Factor: 1 Analyst: TLS
 Injection Volume: 1 (µL) Prep Method: _____
 GPC Cleanup: (Y/N) N pH: _____ Analytical Method: OLMO 4.2
 Prep Batch: 271129 Analytical Batch: 271697 Sulfur Cleanup: (Y/N) N Instrument ID: GCS6A
 Lab File ID: 2040331/SV6034

CONCENTRATION UNITS: ug/L

CAS NO. COMPOUND

RESULT

Q

MDL

RL

| | | | | | |
|------------|---------------------|-------|---|---------|-------|
| 72-54-8 | 4,4'-DDD | 0.100 | U | 0.00010 | 0.100 |
| 72-55-9 | 4,4'-DDE | 0.100 | U | 0.00010 | 0.100 |
| 50-29-3 | 4,4'-DDT | 0.100 | U | 0.00010 | 0.100 |
| 309-00-2 | Aldrin | 0.050 | U | 0.00010 | 0.050 |
| 12674-11-2 | Aroclor-1016 | 1.00 | U | 0.00010 | 1.00 |
| 11104-26-2 | Aroclor-1221 | 2.00 | U | 0.00010 | 2.00 |
| 11141-16-5 | Aroclor-1232 | 1.00 | U | 0.00010 | 1.00 |
| 53468-21-9 | Aroclor-1242 | 1.00 | U | 0.00010 | 1.00 |
| 12672-29-6 | Aroclor-1248 | 1.00 | U | 0.00010 | 1.00 |
| 11097-69-1 | Aroclor-1254 | 1.00 | U | 0.00010 | 1.00 |
| 11096-82-5 | Aroclor-1260 | 1.00 | U | 0.00010 | 1.00 |
| 60-57-1 | Dieldrin | 0.100 | U | 0.00010 | 0.100 |
| 959-98-8 | Endosulfan I | 0.050 | U | 0.00010 | 0.050 |
| 33213-65-9 | Endosulfan II | 0.100 | U | 0.00010 | 0.100 |
| 1031-07-8 | Endosulfan sulfate | 0.100 | U | 0.00010 | 0.100 |
| 72-20-8 | Endrin | 0.100 | U | 0.00010 | 0.100 |
| 7421-93-4 | Endrin aldehyde | 0.100 | U | 0.00010 | 0.100 |
| 53494-70-5 | Endrin ketone | 0.100 | U | 0.00010 | 0.100 |
| 76-44-8 | Heptachlor | 0.050 | U | 0.00010 | 0.050 |
| 1024-57-3 | Heptachlor epoxide | 0.050 | U | 0.00010 | 0.050 |
| 72-43-5 | Methoxychlor | 0.500 | U | 0.00010 | 0.500 |
| 6001-35-2 | Toxaphene | 5.00 | U | 0.00010 | 5.00 |
| 319-84-6 | alpha-BHC | 0.050 | U | 0.00010 | 0.050 |
| 5103-71-9 | alpha-Chlordane | 0.050 | U | 0.00010 | 0.050 |
| 319-85-7 | beta-BHC | 0.050 | U | 0.00010 | 0.050 |
| 319-86-8 | delta-BHC | 0.050 | U | 0.00010 | 0.050 |
| 58-89-9 | gamma-BHC (Lincane) | 0.050 | U | 0.00010 | 0.050 |
| 5103-74-2 | gamma-Chlordane | 0.050 | U | 0.00010 | 0.050 |

1D
ORGANICS ANALYSIS DATA SHEET

Lab Name: GCAL Sample ID: SKGW601009
 Lab Code: LA024 Case No.: _____ Contract: _____
 Matrix: Water SAS No.: _____ SDG No.: 204031909
 Sample wt/vol: 1000 Units: mL Lab Sample ID: 20403220804
 Level: (low/med) _____ Date Collected: 03/17/04 Time: 1115
 % Moisture: _____ decanted: (Y/N) _____ Date Received: 03/19/04
 GC Column: DB-608-30M ID: .53 (mm) Date Extracted: 3/23/04
 Concentrated Extract Volume: 1000 (µL) Date Analyzed: 04/01/04 Time: 1013
 Soil Aliquot Volume: _____ (µL) Dilution Factor: 1 Analyst: TLS
 Injection Volume: 1 (µL) Prep Method: _____
 GPC Cleanup: (Y/N) N pH: _____ Analytical Method: OLMO 4.2
 Prep Batch: 271129 Analytical Batch: 271697 Sulfur Cleanup: (Y/N) N Instrument ID: GCS6A
 Lab File ID: 2040331/SV6035

CONCENTRATION UNITS: ug/L

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|------------|---------------------|--------|---|---------|-------|
| 72-54-8 | 4,4'-DDD | 0.100 | U | 0.00010 | 0.100 |
| 72-55-9 | 4,4'-DDE | 0.100 | U | 0.00010 | 0.100 |
| 50-29-3 | 4,4'-DDT | 0.100 | U | 0.00010 | 0.100 |
| 309-00-2 | Aldrin | 0.050 | U | 0.00010 | 0.050 |
| 12674-11-2 | Aroclor-1016 | 1.00 | U | 0.00010 | 1.00 |
| 11104-28-2 | Aroclor-1221 | 2.00 | U | 0.00010 | 2.00 |
| 11141-16-5 | Aroclor-1232 | 1.00 | U | 0.00010 | 1.00 |
| 53469-21-9 | Aroclor-1242 | 1.00 | U | 0.00010 | 1.00 |
| 12672-29-6 | Aroclor-1248 | 1.00 | U | 0.00010 | 1.00 |
| 11097-69-1 | Aroclor-1254 | 1.00 | U | 0.00010 | 1.00 |
| 11096-82-5 | Aroclor-1260 | 1.00 | U | 0.00010 | 1.00 |
| 60-57-1 | Dieldrin | 0.100 | U | 0.00010 | 0.100 |
| 959-98-8 | Endosulfan I | 0.050 | U | 0.00010 | 0.050 |
| 33213-65-9 | Endosulfan II | 0.100 | U | 0.00010 | 0.100 |
| 1031-07-8 | Endosulfan sulfate | 0.100 | U | 0.00010 | 0.100 |
| 72-20-8 | Endrin | 0.100 | U | 0.00010 | 0.100 |
| 7421-93-4 | Endrin aldehyde | 0.100 | U | 0.00010 | 0.100 |
| 53494-70-5 | Endrin sulfate | 0.100 | U | 0.00010 | 0.100 |
| 76-44-8 | Heptachlor | 0.050 | U | 0.00010 | 0.050 |
| 1024-57-3 | Heptachlor epoxide | 0.050 | U | 0.00010 | 0.050 |
| 72-43-5 | Methoxychlor | 0.500 | U | 0.00010 | 0.500 |
| 8001-35-2 | Toxaphene | 5.00 | U | 0.00010 | 5.00 |
| 319-84-6 | alpha-BHC | 0.050 | U | 0.00010 | 0.050 |
| 5103-71-9 | alpha-Chlordane | 0.050 | U | 0.00010 | 0.050 |
| 319-85-7 | beta-BHC | 0.050 | U | 0.00010 | 0.050 |
| 319-86-8 | delta-BHC | 0.050 | U | 0.00010 | 0.050 |
| 58-89-9 | gamma-BHC (Lindane) | 0.050 | U | 0.00010 | 0.050 |
| 5103-74-2 | gamma-Chlordane | 0.050 | U | 0.00010 | 0.050 |

ORGANICS ANALYSIS DATA SHEET

Lab Name: GCAL Sample ID: SKGW62A1009

Lab Code: LA024 Case No.: _____ Contract: _____

Matrix: Water SAS No.: _____ SDG No.: 204031909

Sample w/vol: 1000 Units: mL Lab Sample ID: 20403220805

Level: (low/med) _____ Date Collected: 03/17/04 Time: 1150

% Moisture: _____ decanted: (Y/N) _____ Date Received: 03/19/04

GC Column: DB-608-30M ID: 53 (mm) Date Extracted: 3/23/04

Concentrated Extract Volume: 1000 (µL) Date Analyzed: 04/01/04 Time: 1041

Soil Aliquot Volume: _____ (µL) Dilution Factor: 1 Analyst: TLS

Injection Volume: 1 (µL) Prep Method: _____

GPC Cleanup: (Y/N) N pH: _____ Analytical Method: OLMO 4.2

Prep Batch: 271129 Analytical Batch: 271697 Sulfur Cleanup: (Y/N) N Instrument ID: GCS6A

Lab File ID: 2040331/SV6036

CONCENTRATION UNITS: ug/L

CAS NO. COMPOUND

RESULT

Q

MDL

RL

| | | | | | |
|------------|---------------------|-------|---|---------|-------|
| 72-54-8 | 4,4'-DDD | 0.100 | U | 0.00010 | 0.100 |
| 72-55-9 | 4,4'-DDE | 0.100 | U | 0.00010 | 0.100 |
| 50-29-3 | 4,4'-DDT | 0.100 | U | 0.00010 | 0.100 |
| 309-00-2 | Aldrin | 0.050 | U | 0.00010 | 0.050 |
| 12674-11-2 | Aroclor-1016 | 1.00 | U | 0.00010 | 1.00 |
| 11104-28-2 | Aroclor-1221 | 2.00 | U | 0.00010 | 2.00 |
| 11141-16-5 | Aroclor-1232 | 1.00 | U | 0.00010 | 1.00 |
| 53489-21-9 | Aroclor-1242 | 1.00 | U | 0.00010 | 1.00 |
| 12672-29-6 | Aroclor-1248 | 1.00 | U | 0.00010 | 1.00 |
| 11097-69-1 | Aroclor-1254 | 1.00 | U | 0.00010 | 1.00 |
| 11096-82-5 | Aroclor-1260 | 1.00 | U | 0.00010 | 1.00 |
| 60-57-1 | Dieldrin | 0.100 | U | 0.00010 | 0.100 |
| 958-98-8 | Endosulfan I | 0.050 | U | 0.00010 | 0.050 |
| 33213-85-9 | Endosulfan II | 0.100 | U | 0.00010 | 0.100 |
| 1031-07-8 | Endosulfan sulfate | 0.100 | U | 0.00010 | 0.100 |
| 72-20-8 | Endrin | 0.100 | U | 0.00010 | 0.100 |
| 7421-93-4 | Endrin aldehyde | 0.100 | U | 0.00010 | 0.100 |
| 53494-70-5 | Endrin ketone | 0.100 | U | 0.00010 | 0.100 |
| 76-44-8 | Heptachlor | 0.050 | U | 0.00010 | 0.050 |
| 1024-57-3 | Heptachlor epoxide | 0.050 | U | 0.00010 | 0.050 |
| 72-43-5 | Methoxychlor | 0.500 | U | 0.00010 | 0.500 |
| 8001-35-2 | Toxaphene | 5.00 | U | 0.00010 | 5.00 |
| 319-84-6 | alpha-BHC | 0.050 | U | 0.00010 | 0.050 |
| 5103-71-9 | alpha-Chlordane | 0.050 | U | 0.00010 | 0.050 |
| 319-85-7 | beta-BHC | 0.050 | U | 0.00010 | 0.050 |
| 319-86-8 | delta-BHC | 0.050 | U | 0.00010 | 0.050 |
| 58-89-9 | gamma-BHC (Lindane) | 0.050 | U | 0.00010 | 0.050 |
| 5103-74-2 | gamma-Chlordane | 0.050 | U | 0.00010 | 0.050 |

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ORGANICS ANALYSIS DATA SHEET

Lab Name: GCAL Sample ID: SKGW641009
 Lab Code: LA024 Case No.: _____ Contract: _____
 Matrix: Water SAS No.: _____ SDG No.: 204031909
 Sample wt/vol: 1000 Units: mL Lab Sample ID: 20403220806
 Level: (low/med) _____ Date Collected: 03/17/04 Time: 1410
 % Moisture: _____ decanted: (Y/N) _____ Date Received: 03/19/04
 GC Column: DB-608-30M ID: .53 (mm) Date Extracted: 3/23/04
 Concentrated Extract Volume: 1000 (µL) Date Analyzed: 04/01/04 Time: 1109
 Soil Aliquot Volume: _____ (µL) Dilution Factor: 1 Analyst: TLS
 Injection Volume: 1 (µL) Prep Method: _____
 GPC Cleanup: (Y/N) N pH: _____ Analytical Method: OLMO 4.2
 Prep Batch: 271129 Analytical Batch: 271697 Sulfur Cleanup: (Y/N) N Instrument ID: GCS6A
 CONCENTRATION UNITS: ug/L Lab File ID: 2040331/SV6037

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|------------|---------------------|--------|---|---------|-------|
| 72-54-8 | 4,4'-DDD | 0.100 | U | 0.00010 | 0.100 |
| 72-55-9 | 4,4'-DDE | 0.100 | U | 0.00010 | 0.100 |
| 50-29-3 | 4,4'-DDT | 0.100 | U | 0.00010 | 0.100 |
| 309-00-2 | Aldrin | 0.050 | U | 0.00010 | 0.050 |
| 12674-11-2 | Aroclor-1018 | 1.00 | U | 0.00010 | 1.00 |
| 11104-28-2 | Aroclor-1221 | 2.00 | U | 0.00010 | 2.00 |
| 11141-16-5 | Aroclor-1232 | 1.00 | U | 0.00010 | 1.00 |
| 53469-21-9 | Aroclor-1242 | 1.00 | U | 0.00010 | 1.00 |
| 12672-29-6 | Aroclor-1248 | 1.00 | U | 0.00010 | 1.00 |
| 11097-69-1 | Aroclor-1254 | 1.00 | U | 0.00010 | 1.00 |
| 11096-82-5 | Aroclor-1260 | 1.00 | U | 0.00010 | 1.00 |
| 60-57-1 | Dieldrin | 0.100 | U | 0.00010 | 0.100 |
| 959-98-8 | Endosulfan I | 0.050 | U | 0.00010 | 0.050 |
| 33213-65-9 | Endosulfan II | 0.100 | U | 0.00010 | 0.100 |
| 1031-07-8 | Endosulfan sulfate | 0.100 | U | 0.00010 | 0.100 |
| 72-20-8 | Endrin | 0.100 | U | 0.00010 | 0.100 |
| 7421-93-4 | Endrin aldehyde | 0.100 | U | 0.00010 | 0.100 |
| 53494-70-5 | Endrin ketone | 0.100 | U | 0.00010 | 0.100 |
| 76-44-8 | Heptachlor | 0.050 | U | 0.00010 | 0.050 |
| 1024-57-3 | Heptachlor epoxide | 0.050 | U | 0.00010 | 0.050 |
| 72-43-5 | Methoxychlor | 0.500 | U | 0.00010 | 0.500 |
| 8001-35-2 | Toxaphene | 5.00 | U | 0.00010 | 5.00 |
| 319-84-6 | alpha-BHC | 0.050 | U | 0.00010 | 0.050 |
| 5103-71-9 | alpha-Chlordane | 0.050 | U | 0.00010 | 0.050 |
| 319-85-7 | beta-BHC | 0.050 | U | 0.00010 | 0.050 |
| 319-86-8 | delta-BHC | 0.050 | U | 0.00010 | 0.050 |
| 58-89-9 | gamma-BHC (Lindane) | 0.050 | U | 0.00010 | 0.050 |
| 5103-74-2 | gamma-Chlordane | 0.050 | U | 0.00010 | 0.050 |

INORGANIC ANALYSIS DATA SHEET

SKGW581009

Lab Name: GCAL Contract: _____
 Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204031909
 Matrix (soil / water): Water Lab Sample ID: 20403190901
 Level: (low / med) _____ Date Received: 03/19/04
 % Solids: _____

Concentration Units (ug/L or mg/kg dry weight): ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7440-36-0 | Antimony | 5.7 | B | | P |
| 7440-38-2 | Arsenic | 11.5 | | | P |
| 7440-39-3 | Barium | 284 | | | P |
| 7440-41-7 | Beryllium | 1.0 | B | | P |
| 7440-43-9 | Cadmium | 1.5 | B | | P |
| 7440-47-3 | Chromium | 28.2 | | | P |
| 7440-50-8 | Copper | 45.7 | | | P |
| 7439-89-6 | Iron | 32700 | | | P |
| 7439-92-1 | Lead | 19.5 | | | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 32.1 | B | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-28-0 | Thallium | 4.1 | B | | P |
| 7440-66-6 | Zinc | 81.0 | | | P |
| 57-12-5 | Cyanide | 0.5 | U | | AS |

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Color Before: LT. BROWN Clarity Before: CLEAR Texture: _____
 Color After: LT. BROWN Clarity After: CLEAR Artifacts: _____
 Comments: _____

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INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKGW58DUP1009

Lab Name: GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204031909
Matrix: (soil / water) Water Lab Sample ID: 20403190902
Level: (low / med) _____ Date Received: 03/19/04
% Solids: _____
Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7440-36-0 | Antimony | 3.7 | B | | P |
| 7440-38-2 | Arsenic | 15.1 | | | P |
| 7440-39-3 | Barium | 326 | | | P |
| 7440-41-7 | Beryllium | 1.3 | B | | P |
| 7440-43-9 | Cadmium | 1.9 | B | | P |
| 7440-47-3 | Chromium | 41.8 | | | P |
| 7440-50-8 | Copper | 55.9 | | | P |
| 7439-89-8 | Iron | 48400 | | | P |
| 7439-92-1 | Lead | 27.2 | | | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 46.4 | | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-66-8 | Zinc | 125 | | | P |
| 57-12-5 | Cyanide | 0.5 | U | | AS |

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Color Before: LT. BROWN Clarity Before: CLEAR Texture: _____
Color After: LT. BROWN Clarity After: CLEAR Artifacts: _____
Comments: _____

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INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKGW631009

Lab Name: GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204031909
Metric (soil / water) Water Lab Sample ID: 20403190903
Level: (low / med) _____ Date Received: 03/19/04
% Solids: _____

Concentration Units (ug/L or mg/kg dry weight): ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7440-38-0 | Antimony | 5.7 | B | | P |
| 7440-38-2 | Arsenic | 17.1 | | | P |
| 7440-39-3 | Barium | 186 | B | | P |
| 7440-41-7 | Beryllium | 2.1 | B | | P |
| 7440-43-9 | Cadmium | 2.5 | B | | P |
| 7440-47-3 | Chromium | 38.2 | | | P |
| 7440-50-8 | Copper | 69.2 | | | P |
| 7439-89-6 | Iron | 63200 | | | P |
| 7439-92-1 | Lead | 41.0 | | | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 58.1 | | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-28-0 | Thallium | 8.5 | B | | P |
| 7440-66-6 | Zinc | 176 | | | P |
| 57-12-5 | Cyanide | 0.5 | U | | AS |

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Color Before: LT. BROWN Clarity Before: CLEAR Texture: _____
Color After: LT. BROWN Clarity After: CLEAR Artifacts: _____
Comments: _____

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INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKGW611009

Lab Name: GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204031909
Matrix: (soil / water) Water Lab Sample ID: 20403190904
Level: (low / med) _____ Date Received: 03/19/04
% Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7440-36-0 | Antimony | 4.8 | B | | P |
| 7440-38-2 | Arsenic | 2.9 | U | | P |
| 7440-39-3 | Barium | 44.1 | B | | P |
| 7440-41-7 | Beryllium | 0.2 | B | | P |
| 7440-43-9 | Cadmium | 0.3 | B | | P |
| 7440-47-3 | Chromium | 1.9 | B | | P |
| 7440-50-8 | Copper | 22.2 | B | | P |
| 7439-89-6 | Iron | 2430 | | | P |
| 7439-92-1 | Lead | 22.1 | | | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 4.3 | B | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-66-6 | Zinc | 7.3 | B | | P |
| 57-12-5 | Cyanide | 0.5 | U | | AS |

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Color Before: COLORLESS Clarity Before: CLEAR Texture: _____
Color After: COLORLESS Clarity After: CLEAR Artifacts: _____
Comments: _____

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INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKGW61DUP1009

Lab Name: GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204031909
Matrix (soil / water) Water Lab Sample ID 20403190907
Level: (low / med) _____ Date Received 03/19/04
% Solids: _____
Concentration Units (ug/L or mg/kg dry weight): ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7440-38-0 | Antimony | 5.5 | B | | P |
| 7440-38-2 | Arsenic | 3.7 | B | | P |
| 7440-39-3 | Barium | 45.4 | B | | P |
| 7440-41-7 | Beryllium | 0.2 | B | | P |
| 7440-43-0 | Cadmium | 0.3 | B | | P |
| 7440-47-3 | Chromium | 1.6 | B | | P |
| 7440-50-8 | Copper | 18.6 | B | | P |
| 7439-89-6 | Iron | 2490 | | | P |
| 7439-82-1 | Lead | 22.5 | | | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 3.7 | B | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-66-6 | Zinc | 7.0 | B | | P |
| 57-12-5 | Cyanide | 0.8 | B | | AS |

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Color Before: COLORLESS Clarity Before: CLEAR Texture: _____
Color After: COLORLESS Clarity After: CLEAR Artifacts: _____
Comments: _____

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INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKGWFB1009

Lab Name: GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204031909
Matrix: (soil / water) Water Lab Sample ID: 20403190908
Level: (low / med) _____ Date Received: 03/19/04
% Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7440-36-0 | Antimony | 3.7 | U | | P |
| 7440-38-2 | Arsenic | 2.9 | U | | P |
| 7440-39-3 | Barium | 0.4 | B | | P |
| 7440-41-7 | Beryllium | 0.1 | U | | P |
| 7440-43-9 | Cadmium | 0.2 | U | | P |
| 7440-47-3 | Chromium | 0.8 | U | | P |
| 7440-50-8 | Copper | 1.2 | U | | P |
| 7439-89-6 | Iron | 14.1 | U | | P |
| 7439-92-1 | Lead | 1.5 | U | | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 0.9 | B | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-28-0 | Thallium | 4.6 | B | | P |
| 7440-66-6 | Zinc | 3.2 | B | | P |
| 57-12-5 | Cyanide | 0.5 | U | | AS |

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Color Before: COLORLESS Clarity Before: CLEAR Texture: _____
Color After: COLORLESS Clarity After: CLEAR Artifacts: _____
Comments: _____

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EPA SAMPLE NO.

INORGANIC ANALYSIS DATA SHEET

SKGW581009(DISS)

Lab Name: GCALContract Lab Code: LA024Case No.: SAS No.: SDG No.: 204031909Metric (soil / water) WaterLab Sample ID: 20403190910Level: (low / med) Date Received: 03/19/04% Solids: Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7440-38-0 | Antimony | 3.7 | U | | P |
| 7440-38-2 | Arsenic | 3.1 | B | | P |
| 7440-39-3 | Barium | 156 | B | | P |
| 7440-41-7 | Beryllium | 0.1 | U | | P |
| 7440-43-8 | Cadmium | 0.2 | U | | P |
| 7440-47-3 | Chromium | 1.5 | B | | P |
| 7440-50-8 | Copper | 2.9 | B | | P |
| 7439-89-6 | Iron | 209 | | | P |
| 7439-92-1 | Lead | 1.5 | U | | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 2.6 | B | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-66-6 | Zinc | 0.6 | U | | P |

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Color Before: COLORLESSClarity Before: CLEARTexture: Color After: COLORLESSClarity After: CLEARArtifacts: Comments:

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INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKGW58DUP1009(DISS)

Lab Name: GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204031909
Matrix: (soil / water) Water Lab Sample ID: 20403190911
Level: (low / med) _____ Date Received: 03/19/04
% Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7440-36-0 | Antimony | 3.7 | U | | P |
| 7440-38-2 | Arsenic | 2.9 | U | | P |
| 7440-39-3 | Barium | 164 | B | | P |
| 7440-41-7 | Beryllium | 0.1 | U | | P |
| 7440-43-9 | Cadmium | 0.2 | U | | P |
| 7440-47-3 | Chromium | 1.3 | B | | P |
| 7440-50-8 | Copper | 1.2 | U | | P |
| 7439-89-6 | Iron | 201 | | | P |
| 7439-92-1 | Lead | 1.5 | U | | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 2.0 | B | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-66-6 | Zinc | 0.6 | U | | P |

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Color Before: COLORLESS Clarity Before: CLEAR Texture: _____
Color After: COLORLESS Clarity After: CLEAR Artifacts: _____
Comments: _____

U.S. EPA - CLP
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKGW631009(DISS)

Lab Name: GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204031909
Matrix (soil / water): Water Lab Sample ID: 20403190912
Level: (low / med) _____ Date Received: 03/19/04
% Solids: _____

Concentration Units (ug/L or mg/kg dry weight): ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7440-36-0 | Antimony | 3.8 | B | | P |
| 7440-38-2 | Arsenic | 2.9 | U | | P |
| 7440-39-3 | Barium | 20.1 | B | | P |
| 7440-41-7 | Beryllium | 0.2 | B | | P |
| 7440-43-8 | Cadmium | 0.2 | U | | P |
| 7440-47-3 | Chromium | 1.8 | B | | P |
| 7440-50-8 | Copper | 2.0 | B | | P |
| 7439-89-6 | Iron | 21.4 | B | | P |
| 7439-92-1 | Lead | 1.5 | U | | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 3.2 | B | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-28-0 | Thallium | 4.6 | B | | P |
| 7440-66-6 | Zinc | 0.6 | U | | P |

43

5/25/04
m/s

Color Before: COLORLESS Clarity Before: CLEAR Texture: _____
Color After: COLORLESS Clarity After: CLEAR Artifacts: _____
Comments: _____

U.S. EPA - CLP
1
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKGW611009(DISS)

Lab Name: GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204031909
Matrix: (soil / water) Water Lab Sample ID: 20403190913
Level: (low / med) _____ Date Received: 03/19/04
% Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7440-36-0 | Antimony | 4.5 | B | | P |
| 7440-38-2 | Arsenic | 2.9 | U | | P |
| 7440-39-3 | Barium | 39.4 | B | | P |
| 7440-41-7 | Beryllium | 0.1 | B | | P |
| 7440-43-9 | Cadmium | 0.3 | B | | P |
| 7440-47-3 | Chromium | 1.1 | B | | P |
| 7440-50-8 | Copper | 8.0 | B | | P |
| 7439-89-6 | Iron | 187 | | | P |
| 7439-92-1 | Lead | 1.5 | U | | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 4.2 | B | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-66-6 | Zinc | 0.6 | U | | P |

US

5/28/04
me

Color Before: COLORLESS Clarity Before: CLEAR Texture: _____
Color After: COLORLESS Clarity After: CLEAR Artifacts: _____
Comments: _____

U.S. EPA - CLP

EPA SAMPLE NO.

INORGANIC ANALYSIS DATA SHEET

SKGW61DUP1009 (DISS)

Lab Name: GCAL

Contract: _____

Lab Code: LA024

Case No.: _____

SAS No.: _____

SDG No.: 204031909Matrix (soil / water) WaterLab Sample ID 20403190915

Level: (low / med) _____

Date Received: 03/19/04

% Solids: _____

Concentration Units (ug/L or mg/kg dry weight): ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7440-38-0 | Antimony | 3.7 | U | | P |
| 7440-38-2 | Arsenic | 2.9 | U | | P |
| 7440-39-3 | Barium | 38.9 | B | | P |
| 7440-41-7 | Beryllium | 0.1 | B | | P |
| 7440-43-9 | Cadmium | 0.2 | U | | P |
| 7440-47-3 | Chromium | 1.3 | B | | P |
| 7440-50-8 | Copper | 5.2 | B | | P |
| 7439-89-6 | Iron | 195 | | | P |
| 7439-92-1 | Lead | 1.5 | U | | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 3.6 | B | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-28-0 | Thallium | 5.6 | B | | P |
| 7440-66-6 | Zinc | 0.6 | U | | P |

u/s

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5/25/04
mizColor Before: COLORLESSClarity Before: CLEAR

Texture: _____

Color After: COLORLESSClarity After: CLEAR

Artifacts: _____

Comments: _____

U.S. EPA - CLP
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INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKGWFB1009 (DISS)

Lab Name: GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204031909
Matrix: (soil / water) Water Lab Sample ID: 20403190916
Level: (low / med) _____ Date Received: 03/19/04
% Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7440-36-0 | Antimony | 3.7 | B | | P |
| 7440-38-2 | Arsenic | 2.9 | U | | P |
| 7440-39-3 | Barium | 0.3 | U | | P |
| 7440-41-7 | Beryllium | 0.1 | U | | P |
| 7440-43-9 | Cadmium | 0.2 | U | | P |
| 7440-47-3 | Chromium | 0.8 | U | | P |
| 7440-50-8 | Copper | 1.2 | U | | P |
| 7439-89-6 | Iron | 14.1 | U | | P |
| 7439-92-1 | Lead | 1.5 | U | | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 0.7 | U | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-28-0 | Thallium | 7.5 | B | | P |
| 7440-66-6 | Zinc | 0.6 | U | | P |

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Color Before: COLORLESS Clarity Before: CLEAR Texture: _____
Color After: COLORLESS Clarity After: CLEAR Artifacts: _____
Comments:

U.S. EPA - CLP
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKGW06R1009

Lab Name: GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204031909
Matrix (soil / water): Water Lab Sample ID: 20403220801
Level: (low / med) _____ Date Received: 03/19/04
% Solids: _____
Concentration Units (ug/L or mg/kg dry weight): ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7440-38-0 | Antimony | 5.5 | B | | P |
| 7440-38-2 | Arsenic | 12.4 | | | P |
| 7440-39-3 | Barium | 440 | | | P |
| 7440-41-7 | Beryllium | 1.1 | B | | P |
| 7440-43-0 | Cadmium | 1.0 | B | | P |
| 7440-47-3 | Chromium | 16.9 | | | P |
| 7440-50-8 | Copper | 39.3 | | | P |
| 7439-89-6 | Iron | 25300 | | | P |
| 7439-92-1 | Lead | 23.9 | | | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 23.4 | B | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-86-6 | Zinc | 72.9 | | | P |
| 57-12-5 | Cyanide | 1.0 | B | | AS |

45

5/25/04
mu

Color Before: LT. YELLOW Clarity Before: CLEAR Texture: _____
Color After: LT. YELLOW Clarity After: CLEAR Artifacts: _____
Comments: _____

U.S. EPA - CLP
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INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKGW07R1009

Lab Name: GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204031909
Matrix: (soil / water) Water Lab Sample ID: 20403220802
Level: (low / med) _____ Date Received: 03/19/04
% Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7440-36-0 | Antimony | 6.6 | B | | P |
| 7440-38-2 | Arsenic | 6.9 | B | | P |
| 7440-39-3 | Barium | 484 | | | P |
| 7440-41-7 | Beryllium | 0.8 | B | | P |
| 7440-43-9 | Cadmium | 0.9 | B | | P |
| 7440-47-3 | Chromium | 12.9 | | | P |
| 7440-50-8 | Copper | 35.5 | | | P |
| 7439-89-6 | Iron | 20200 | | | P |
| 7439-92-1 | Lead | 9.2 | | | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 17.8 | B | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-66-6 | Zinc | 51.2 | | | P |
| 57-12-5 | Cyanide | 1.5 | B | | AS |

WS

5/25/04
me

Color Before: LT.YELLOW Clarity Before: CLEAR Texture: _____
Color After: LT.YELLOW Clarity After: CLEAR Artifacts: _____
Comments: _____

U.S. EPA - CLP
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKGW591009

Lab Name: GCAL Contract: _____
Lab Code: LA024 Case No. _____ SAS No.: _____ SDG No.: 204031909
Matrix: (soil / water) Water Lab Sample ID: 20403220803
Level: (low / med) _____ Date Received: 03/19/04
% Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7440-36-0 | Antimony | 4.7 | B | | P |
| 7440-38-2 | Arsenic | 2.9 | U | | P |
| 7440-39-3 | Barium | 55.0 | B | | P |
| 7440-41-7 | Beryllium | 0.2 | B | | P |
| 7440-43-8 | Cadmium | 0.2 | U | | P |
| 7440-47-3 | Chromium | 5.5 | B | | P |
| 7440-50-8 | Copper | 10.1 | B | | P |
| 7439-89-6 | Iron | 3020 | | | P |
| 7439-92-1 | Lead | 1.5 | U | | P |
| 7439-97-8 | Mercury | 0.1 | L | | AV |
| 7440-02-0 | Nickel | 6.7 | B | | P |
| 7782-49-2 | Selenium | 4.4 | L | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-28-0 | Thallium | 2.6 | L | | P |
| 7440-66-6 | Zinc | 7.9 | B | | P |
| 57-12-5 | Cyanide | 1.0 | B | | AS |

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5/28/04
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Color Before: COLORLESS Clarity Before: CLEAR Texture: _____
Color After: COLORLESS Clarity After: CLEAR Artifacts: _____
Comments: _____

U.S. EPA - CLP
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INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKGW601009

Lab Name: GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204031909
Matrix: (soil / water) Water Lab Sample ID: 20403220804
Level: (low / med) _____ Date Received: 03/19/04
% Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7440-36-0 | Antimony | 9.7 | B | | P |
| 7440-38-2 | Arsenic | 17.0 | | | P |
| 7440-39-3 | Barium | 129 | B | | P |
| 7440-41-7 | Beryllium | 2.5 | B | | P |
| 7440-43-9 | Cadmium | 2.8 | B | | P |
| 7440-47-3 | Chromium | 59.6 | | | P |
| 7440-50-8 | Copper | 54.5 | | | P |
| 7439-89-8 | Iron | 74200 | | | P |
| 7439-92-1 | Lead | 40.4 | | | P |
| 7439-97-6 | Mercury | 0.1 | B | | AV |
| 7440-02-0 | Nickel | 67.3 | | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-28-0 | Thallium | 11.0 | | | P |
| 7440-66-6 | Zinc | 180 | | | P |

US

5/28/04
m

Color Before: DK. BROWN Clarity Before: CLEAR Texture: _____
Color After: DK. BROWN Clarity After: CLEAR Artifacts: _____
Comments: _____

U.S. EPA - CLP

EPA SAMPLE NO.

INORGANIC ANALYSIS DATA SHEET

SKGW62A1009

Lab Name: GCAL

Contract: _____

Lab Code: LA024

Case No.: _____

SAS No.: _____

SDG No.: 204031909Matrix (soil / water) WaterLab Sample ID: 20403220805

Level: (low / med) _____

Date Received 03/19/04

% Solids: _____

Concentration Units (ug/L or mg/kg dry weight): ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7440-38-0 | Antimony | 6.1 | B | | P |
| 7440-38-2 | Arsenic | 8.3 | B | | P |
| 7440-39-3 | Barium | 361 | | | P |
| 7440-41-7 | Beryllium | 1.1 | B | | P |
| 7440-43-8 | Cadmium | 1.6 | B | | P |
| 7440-47-3 | Chromium | 29.6 | | | P |
| 7440-50-8 | Copper | 42.7 | | | P |
| 7439-89-8 | Iron | 35000 | | | P |
| 7439-92-1 | Lead | 39.5 | | | P |
| 7439-97-8 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 35.4 | B | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-28-0 | Thallium | 6.9 | B | | P |
| 7440-66-6 | Zinc | 101 | | | P |
| 57-12-5 | Cyanide | 1.0 | B | | AS |

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5/28/04
micColor Before: LT. BROWNClarity Before: CLEAR

Texture: _____

Color After: LT. BROWNClarity After: CLEAR

Artifacts: _____

Comments: _____

U.S. EPA - CLP
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INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKGW641009

Lab Name: GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204031909
Matrix: (soil / water) Water Lab Sample ID: 20403220806
Level: (low / med) _____ Date Received: 03/19/04
% Solids: _____
Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7440-36-0 | Antimony | 4.9 | B | | P |
| 7440-38-2 | Arsenic | 2.9 | U | | P |
| 7440-39-3 | Barium | 37.1 | B | | P |
| 7440-41-7 | Beryllium | 0.3 | B | | P |
| 7440-43-9 | Cadmium | 0.2 | U | | P |
| 7440-47-3 | Chromium | 7.0 | B | | P |
| 7440-50-8 | Copper | 11.3 | B | | P |
| 7439-89-6 | Iron | 7520 | | | P |
| 7439-92-1 | Lead | 1.5 | U | | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 16.4 | B | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-66-6 | Zinc | 13.6 | B | | P |
| 57-12-5 | Cyanide | 1.3 | B | | AS |

43

5

just
me

Color Before: COLORLESS

Color Before: CLEAR

Texture: _____

Color After: COLORLESS

Color After: CLEAR

Artifacts: _____

Comments:

U.S. EPA - C.P.
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKGW06R1009 (DISS)

Lab Name: GCAL Contract: _____
Lab Code: LA024 Case No: _____ SAS No: _____ SDG No.: 204031909
Matrix: (soil / water) Water Lab Sample ID: 20403220808
Level: (low / med) _____ Date Received: 03/19/04
% Solids: _____

Concentration Units (ug/L or mg/kg dry weight): ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7440-38-0 | Antimony | 3.7 | U | | P |
| 7440-38-2 | Arsenic | 2.9 | U | | P |
| 7440-39-3 | Barium | 266 | | | P |
| 7440-41-7 | Beryllium | 0.1 | U | | P |
| 7440-43-8 | Cadmium | 0.2 | U | | P |
| 7440-47-3 | Chromium | 1.2 | B | | P |
| 7440-50-8 | Copper | 1.2 | U | | P |
| 7439-89-6 | Iron | 22.0 | B | | P |
| 7439-92-1 | Lead | 1.5 | U | | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 1.7 | B | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-66-6 | Zinc | 0.6 | U | | P |

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5/28/04
m.e.

Color Before: COLORLESS Clarity Before: CLEAR Texture: _____
Color After: COLORLESS Clarity After: CLEAR Artifacts: _____
Comments: _____

U.S. EPA - CLP
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INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKGW07R1009 (DISS)

Lab Name: GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204031909
Matrix: (soil / water) Water Lab Sample ID: 20403220809
Level: (low / med) _____ Date Received: 03/19/04
% Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7440-36-0 | Antimony | 3.7 | U | | P |
| 7440-38-2 | Arsenic | 2.9 | U | | P |
| 7440-39-3 | Barium | 113 | B | | P |
| 7440-41-7 | Beryllium | 0.1 | B | | P |
| 7440-43-9 | Cadmium | 0.2 | U | | P |
| 7440-47-3 | Chromium | 1.1 | B | | P |
| 7440-50-8 | Copper | 1.2 | U | | P |
| 7439-89-6 | Iron | 32.9 | B | | P |
| 7439-92-1 | Lead | 1.5 | U | | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 2.2 | B | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-66-6 | Zinc | 0.6 | U | | P |

WS

5/20/04
M

Color Before: COLORLESS Clarity Before: CLEAR Texture: _____
Color After: COLORLESS Clarity After: CLEAR Artifacts: _____
Comments:

U.S. EPA - CLP
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INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKGW591009 (DISS)

Lab Name: GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204031909
Metric (soil / water) Water Lab Sample ID: 20403220810
Level: (low / med) _____ Date Received: 03/19/04
% Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7440-38-0 | Antimony | 3.7 | U | | P |
| 7440-38-2 | Arsenic | 2.9 | U | | P |
| 7440-39-3 | Barium | 21.8 | E | | P |
| 7440-41-7 | Beryllium | 0.1 | B | | P |
| 7440-43-8 | Cadmium | 0.2 | L | | P |
| 7440-47-3 | Chromium | 1.8 | B | | P |
| 7440-50-8 | Copper | 2.1 | B | | P |
| 7439-89-6 | Iron | 28.8 | B | | P |
| 7439-92-1 | Lead | 1.5 | U | | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 2.6 | B | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | L | | P |
| 7440-28-0 | Thallium | 3.1 | B | | P |
| 7440-66-6 | Zinc | 3.1 | B | | P |

US

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4/5/04
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Color Before: COLORLESS Clarity Before: CLEAR Texture: _____
Color After: COLORLESS Clarity After: CLEAR Artifacts: _____
Comments: _____

U.S. EPA - CLP
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INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKGW601009 (DISS)

Lab Name: GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204031909
Matrix: (soil / water) Water Lab Sample ID: 20403220811
Level: (low / med) _____ Date Received: 03/19/04
% Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7440-36-0 | Antimony | 5.1 | B | | P |
| 7440-38-2 | Arsenic | 2.9 | U | | P |
| 7440-39-3 | Barium | 27.1 | B | | P |
| 7440-41-7 | Beryllium | 0.2 | B | | P |
| 7440-43-9 | Cadmium | 0.2 | U | | P |
| 7440-47-3 | Chromium | 2.6 | B | | P |
| 7440-50-8 | Copper | 4.0 | B | | P |
| 7439-89-6 | Iron | 14.1 | U | | P |
| 7439-92-1 | Lead | 1.5 | U | | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 2.4 | B | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-66-6 | Zinc | 0.6 | U | | P |

WS

24/04
m+

Color Before: COLORLESS

Clarity Before: CLEAR

Texture: _____

Color After: COLORLESS

Clarity After: CLEAR

Artifacts: _____

Comments:

U.S. EPA - CLP

EPA SAMPLE NO.

INORGANIC ANALYSIS DATA SHEET

SKGW62A1009 (DISS)

Lab Name: GCAL

Contract: _____

Lab Code: LA024

Case No.: _____

SAS No.: _____

SDG No.: 204031909Matrix: (soil / water) WaterLab Sample ID: 20403220812

Level: (low / med) _____

Date Received: 03/19/04

% Solids: _____

Concentration Units (ug/L or mg/kg dry weight): ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7440-38-0 | Antimony | 3.7 | U | | P |
| 7440-38-2 | Arsenic | 2.9 | U | | P |
| 7440-39-3 | Barium | 111 | B | | P |
| 7440-41-7 | Beryllium | 0.1 | U | | P |
| 7440-43-8 | Cadmium | 0.2 | U | | P |
| 7440-47-3 | Chromium | 2.1 | B | | P |
| 7440-50-8 | Copper | 1.2 | U | | P |
| 7439-89-6 | Iron | 14.1 | U | | P |
| 7439-92-1 | Lead | 1.5 | U | | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 1.8 | B | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-66-6 | Zinc | 0.6 | U | | P |

US

Jed
prColor Before: COLORLESSClarity Before: CLEAR

Texture: _____

Color After: COLORLESSClarity After: CLEAR

Artifacts: _____

Comments: _____

U.S. EPA - CLP
1
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKGW641009 (DISS)

Lab Name: GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204031909
Matrix: (soil / water) Water Lab Sample ID: 20403220813
Level: (low / med) _____ Date Received: 03/19/04
% Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7440-38-0 | Antimony | 3.7 | U | | P |
| 7440-38-2 | Arsenic | 2.9 | U | | P |
| 7440-39-3 | Barium | 28.3 | B | | P |
| 7440-41-7 | Beryllium | 0.1 | B | | P |
| 7440-43-9 | Cadmium | 0.2 | U | | P |
| 7440-47-3 | Chromium | 1.7 | B | | P |
| 7440-50-8 | Copper | 1.2 | U | | P |
| 7439-89-6 | Iron | 14.1 | U | | P |
| 7439-92-1 | Lead | 1.5 | U | | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 7.8 | B | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-66-6 | Zinc | 0.6 | U | | P |

US

5/25/04
ms

Color Before: COLORLESS Clarity Before: CLEAR Texture: _____
Color After: COLORLESS Clarity After: CLEAR Artifacts: _____
Comments: _____

CHAIN OF CUSTODY RECORD

Lab use only

use only
Earth Tech
Client Name

4342

Client #

204031904

Workorder #

4/2/04

Due Date

[illegible]

Turn Around Time: ☐ 24-48 hrs. ☐ 3 days ☐ 1 week ☒ Standard ☐ Other

Relinquished by: (Signature)

2-14 14-2-22

Relinquished by (Signature)

Notoulished by: (Signature)

Redox H 2025

Relinquished by: (Signature)

Received by: (Signature)

FILED EX

Booked by: (Signature)

Received by: (Signature)

(Krusch &

Received by: (Signature)

Date:

772

5/7

Date: 3/1

311

Date:

Time:

Title.

5

Time: 02

090

Time:

Note:

By submitting these samples, you agree to the terms and conditions contained in our most recent schedule of services.

2011

3.76 : 3.89

CHAIN OF CUSTODY RECORD

Lab use only

Client Name

Client #

Workorder #

Due Date

Report to:

Client:

Address:

Contact:

Phone:

Fax:

Bill to:

Client:

Address:

Contact:

Phone:**Fax:**

Analytical Requests & Method

Lab use only:

Custody Seal

used ☒ yes ☐ no

in tact ☐ ~~yes~~ ☐ no

Temperature °C

P.O. Number

Project Name/Number

Sampled By:

Lab ID

[illegible]

Turn Around Time: ☐ 24-48 hrs. ☐ 3 days ☐ 1 week ☒ Standard ☐ Other

Relinquished by: (Signature)

Received by: (Signature)

Date:

Time:

Note:

Relinquished by: (Signature)

Received by: (Signature)

Date:

Time:

Relinquished by: (Signature)

Received by: (Signature)

Date:

Time:

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WHITE CLIENT FINAL REPORT — CANARY LABORATORY — PINK CLIENT

AL-06 11/98



GULF COAST ANALYTICAL LABORATORIES, INC.
7979 GSR Avenue, Baton Rouge, Louisiana 70820-7402
Phone 225.789.4900 • Fax 225.787.5717

CHAIN OF CUSTODY RECORD

Lab use only

Earth Tech

Client Name

4342

Client #

204031929

Workorder #

4/6/04

Due Date

Report to:

Client: Earth Tech
Address: 211 W. 1st Street
W. 1st Street
Contact: 1234 5678
Phone: 554 442-2706
Fax: 554 442-2711

Bill to:

Client: IMC
Address: IMC
Contact: IMC
Phone: IMC
Fax: IMC

Analytical Requests & Method

Lab use only:

Custody Seal

used ☒ yes ☐ no

in tact ☒ yes ☐ no

Temperature °C

3

P.O. Number

5429

Project Name/Number

Shiner Landfill IG to 2004

Sampled By:

1234 5678 / 1234 5678

| Matrx | Date | Time (2400) | Sample Description | Preservatives | No. Containers |
|-------|---------|-------------|------------------------------|---------------|----------------|
| W | 3/10/04 | 1430 | X Shiner Landfill IG to 2004 | Various | 7 |
| W | 3/10/04 | 1610 | X Shiner Landfill IG to 2004 | Various | 7 |

| | | | | | | |
|-----------|----------------|------------|------|------|------------------|---------|
| Volatiles | Semi-Volatiles | Pesticides | PCBs | PAHs | Dissolved Metals | Cyanide |
| X | X | X | X | X | X | X |
| X | X | X | X | X | X | X |

Remarks:

Refer to Table 7 (TCL) and Table 8 (TCL) of the Final OSM Plan for the list of analytes

Turn Around Time: ☐ 24-48 hrs. ☐ 3 days ☐ 1 week ☒ Standard ☐ Other

Relinquished by: (Signature)

Relinquished by: (Signature)

Relinquished by: (Signature)

Received by: (Signature)

Received by: (Signature)

Received by: (Signature)

Date:

Date:

Date:

Time:

Time:

Time:

Note:

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CHAIN OF CUSTODY RECORD

Use only
Earth Touch

Client #

704031909

4/2/14

WHITE: CLIENT FINAL REPORT — CANARY: LABORATORY — PINK: CLIENT



GULF COAST ANALYTICAL LABORATORIES, INC.
7979 GBRI Avenue, Baton Rouge, Louisiana 70820-7402
Phone 225.769.4900 • Fax 225.767.5717

CHAIN OF CUSTODY RECORD

Lab use only

Earth Tech

Client Name

4342

Client #

704031902⁴
~~704032203~~

Workorder #

4/2/04

Due Date

Report to:

Client: Earth Tech
Address: 202 Vine Street
Wilder, LA 71076
Contact: P. + 1234567
Phone: 554 442-2311
Fax: 554 442-2311

Client:
Address:
Contact:
Phone:
Fax:

Bill to:

Analytical Requests & Method

Lab use only:

Custody Seal

used ☒ yes ☐ no

In tact ☒ yes ☐ no

Temperature °C 4

P.O. Number

SAZ 60

Project Name/Number

Shinner landfill - 10th. 2004

Sampled By

P. Higgins/C. Cox

| Matrix | Date | Time (2400) | Sample Description | Preservatives | No. Containers | Comments |
|--------|---------|-------------|--------------------|---------------|----------------|--------------|
| W | 3/14/04 | 1458 | SHLW 06R 100g | 1462 | 3 | X |
| W | 3/16/04 | 1552 | SHLW 07R 100g | 1462 | 3 | X |
| W | 3/17/04 | 1000 | SHLW 54 100g | 1462 | 3 | X |
| W | 3/17/04 | 1100 | SHLW 60 100g | 1462 | 3 | X |
| W | 3/17/04 | 1150 | SHLW 62A 100g | 1462 | 3 | X |
| W | 3/17/04 | 1414 | SHLW 64 100g | Various | 5 | X X X X Dimm |
| W | 3/17/04 | 1440 | SHLW 65 100g | Various | 4 | X X X X Dimm |
| W | 3/17/04 | - | Trip Blank | | | |

Remarks:

Refer to Table 7 (TCL) and Table 8 (TAL) in the Final OIM Plan for the list of analytes

3 2 1 2 3 4 5 6 7 14

Turn Around Time: ☐ 24-48 hrs. ☐ 3 days ☐ 1 week ☒ Standard ☐ Other

Relinquished by: (Signature)

Relinquished by: (Signature)

Relinquished by: (Signature)

Received by: (Signature)

Received by: (Signature)

Received by: (Signature)

Date:

Date:

Date:

Time:

Time:

Time:

Note:

By submitting these samples, you agree to the terms and

Due Date

... 51 02 10 39



GULF COAST ANALYTICAL LABORATORIES, INC.
7879 GSRI Avenue, Baton Rouge, Louisiana 70820-7402
Phone 225.769.4800 • Fax 225.767.5717

CHAIN OF CUSTODY RECORD

Lab use only

Earth Tech

Client Name

4342

Client #

30405107 HC
~~304032208~~

Workorder #

4/2/01
Due Date

Report to:

Client: *Earth Tech*
Address: *200 Vine Street*
Wilder, KY 40376
Contact: *P. Higgins*
Phone: *859 442-2300*
Fax: *859 442-2311*

Bill to:

Client: _____
Address: _____
Contact: *(Signature)*
Phone: _____
Fax: _____

Analytical Requests & Method

Lab use only:

Custody Seal

used ☒ yes ☐ no

in tact ☒ yes ☐ no

Temperature °C

P.O. Number

54240

Project Name/Number

Shower 2, well 11 - 104.300.4

Sampled By:

P. Higgins / Chris Cox

| Matrix | Date | Time (2400) | Sample Description | Preservatives | No Containers |
|--------|---------|-------------|--------------------|---------------|---------------|
| W | 3/17/01 | 1054 | Shower 2 100g | Various | 7 |
| W | 3/17/01 | 1155 | Shower 2 60 100g | Various | 6 |

| | | | | | | |
|------------------|-----------------------|-------------|-------------|---------------------|-------------------------|----------------|
| <i>Volatiles</i> | <i>Semi-Volatiles</i> | <i>PSHs</i> | <i>PCBs</i> | <i>total metals</i> | <i>Dissolved Metals</i> | <i>Cyanide</i> |
| X | X | X | X | X | X | X |
| X | X | X | X | X | X | X |

Remarks:

Refer to Table 7 and Table 8, TCC and TAL, & Final OSM Plan for analyte list

Lab ID

3-2-4

Turn Around Time: ☐ 24-48 hrs. ☐ 3 days ☐ 1 week ☒ Standard ☐ Other

Relinquished by: (Signature)

Pat Higgins

Relinquished by: (Signature)

Pat Higgins

Relinquished by: (Signature)

Pat Higgins

Received by: (Signature)

Chris Cox

Received by: (Signature)

Chris Cox

Received by: (Signature)

Chris Cox

Date:

3/17/01

Date:

3/17/01

Date:

3/17/01

Time:

1054

Time:

1155

Time:

1155

Note:

Shower 2 60 100g has low sample volume for Dissolved Metals

By submitting these samples, you agree to the terms and



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Phone 225.769.4900 • Fax 225.767.5717

CHAIN OF CUSTODY RECORD

Lab use only

Earth Tech

Client Name

4342

Client #

304031409
304032208

Workorder #

4/2/08

Due Date

Report to:

Client: Earth Tech
Address: 201 Vine Street
Wilder, KY 41076
Contact: Pat Higgins
Phone: 854 442-2300
Fax: 854 442-2311

Bill to:

Client:
Address:
Contact:
Phone:
Fax:

Analytical Requests & Method

| Volatiles | Semi-Volatiles | Pesticides | PCBs | Total Metals | Dissolved Metals | Cyanide | Class |
|-----------|----------------|------------|------|--------------|------------------|---------|-------------------------------|
| X | X | X | X | X | X | X | 12 |
| X | X | X | X | X | X | X | 13 |
| X | X | X | X | X | X | X | Dimm No sample received |

Lab use only:

Custody Seal
used ☒ yes ☐ no
in tact ☒ yes ☐ no
Temperature °C

P.O. Number 54280 Project Name/Number Skinner Landfill - 1 Qtr. 2004

Sampled By: Pat Higgins / Chris Cox

| Matrix | Date | Time (2400) | C | D | G | B | Sample Description | Preservatives | No Containers |
|--------|---------|-------------|---|---|---|---|--------------------|---------------|---------------|
| W | 3/17/04 | 1150 | | | | | SHGW 62A 100g | Various | 7 |
| W | 3/17/04 | 1410 | | | | | SHGW 64 100g | Various | 5 |
| W | 3/17/04 | 1440 | | | | | SHGW 65 100g | Various | 6 |

Remarks:

Lab ID

3/22

5

6

7

Turn Around Time: ☐ 24-48 hrs. ☐ 3 days ☐ 1 week ☒ Standard ☐ Other

Relinquished by: (Signature)

Pat Higgins

Received by: (Signature)

FED (X)

Date:

3/17/04

Time:

0945

Relinquished by: (Signature)

FED EX # 842586011603

Received by: (Signature)

Chris Cox

Date:

3/19/04

Time:

0945

Note: SHGW 62A - low sample volume for cyanide

By submitting these samples, you agree to the terms and conditions contained in our most recent schedule of services.

DATA VALIDATION REPORT
FOR
SKINNER LANDFILL SITE
EARTH TECH: PROJECT NUMBER 54280
LABORATORY REPORT NUMBER 204030804
PROJECT MANAGER: Ron Rolker
Date: May 6, 2004
Data Validator: Mark Kromis

APPENDIX C LIST OF ACRONYMS

| | |
|--------|--|
| BFB | Bromofluorobenzene |
| CC | Continuing Calibration |
| CCV | Continuing Calibration Verification |
| CCB | Continuing Calibration Blanks |
| CLP | Contract Laboratory Program |
| CRDL | Contract Required Detection Limit |
| DFTPP | Decafluorotriphenylphosphine |
| GC/MS | Gas Chromatograph/Mass Spectrometer |
| IC | Initial Calibration |
| ICB | Initial Calibration Blank |
| IDL | Instrument Detection Limit |
| ICP | Inductively Coupled Plasma |
| ICS | Interference Check Sample |
| ICV | Initial Calibration Verification |
| ILM | Inorganic Analysis Multi-Media Multi-Concentration |
| INDAM | Individual A Mixture |
| INDBM | Individual B Mixture |
| mg/L | milligrams per liter |
| MS/MSD | Matrix Spike/Matrix Spike Duplicate |
| OLC | Organic Analysis Low Concentration |
| OLM | Organic Analysis Multi-Media Multi-Concentration |
| %D | Percent Difference |
| % RSD | Percent Relative Standard Deviation |
| PB | Preparation Blanks |
| QC | Quality Control |
| RF | Response Factor |
| RPD | Relative Percent Difference |
| RRF | Relative Response Factor |
| SDG | Sample Delivery Group |
| SOW | Statement of Work |
| µg/L | micrograms per liter |
| US EPA | United States Environmental Protection Agency |
| VOC | Volatile Organic Compounds |
| VTSR | Validated Time of Sample Receipt |

DATA VALIDATION SUMMARY – SAMPLE DELIVERY GROUP 204030804 INORGANICS

Validation of the inorganics data, as prepared by Gulf Coast Analytical Laboratories (GCAL) for the samples collected from the Skinner Landfill site in March 2004, was conducted by Earth Tech using the National Functional Guidelines for Inorganic Data Review, (US EPA, February, 1994), as appropriate. The results were reported by GCAL under Sample Delivery Group (SDG) 204030804.

| GCAL # | Sample Description |
|--------------|-----------------------|
| 204030804-01 | SKSWD031009 |
| 204030804-02 | SKSWD03D1009 |
| 204030804-03 | SKSWD03MS1009 |
| 204030804-05 | SKSWD03DUP1009 |
| 204030804-06 | SKSWDEB1009 |
| 204030804-08 | SKSWD031009 (DISS) |
| 204030804-09 | SKSWD03D1009 (DISS) |
| 204030804-10 | SKSWD03MS1009 (DISS) |
| 204030804-11 | SKSWD03DUP1009 (DISS) |
| 204030804-12 | SKSWDEB1009 (DISS) |

INTRODUCTION

Analyses of metals were performed according to Contract Laboratory Program (CLP)-Inorganic Analysis Multi-media Multi-concentration ILM04.1 Statement of Work (SOW). Results of the sample analyses are reported by the laboratory as either qualified or unqualified. Unqualified results mean that the reported values maybe used without reservation. Various qualifier codes are used by the laboratory to denote specific information regarding the analytical results.

The data validation process is intended to evaluate the data on a technical basis. The data package also was subjected to an internal laboratory quality review prior to submission to Earth Tech for data validation.

During the validation process, laboratory-qualified and unqualified data are verified against all available supporting documentation. Based on this evaluation, qualifier codes may be added, deleted or modified by the data user. Final results are therefore, either qualified or unqualified. Validator-qualified results are annotated with the following codes in accordance with the Functional Guidelines:

- U The constituent was analyzed for, but was not detected above the level of the associated analytical reporting limit. The associated value is either the sample quantitation limit or the sample detection limit.
- J The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample.

- UJ The analyte was not detected above the reported sample quantitation limit. However, the reported quantitation limit is approximate and may or may not represent the actual limit of quantitation necessary to accurately and precisely measure the analyte in the sample.
- R The sample results are rejected due to serious deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte cannot be verified.

Details of the inorganics data validation findings and conclusions are provided in the following sections of this report:

1. Holding Times
2. Calibration
 - A. Initial Calibration (IC)
 - B. Continuing Calibration (CC)
3. Blanks
4. Inductively Coupled Plasma (ICP) Interference Check Sample
5. Laboratory Control Sample (LCS)
6. Duplicate Analysis
7. Spike Sample Analysis
8. ICP Serial Dilution
9. System Performance
10. Documentation
11. Overall Assessment

1. HOLDING TIMES

All samples for inorganics analyses were analyzed within the 180-day holding time for preserved aqueous samples. Mercury analyses were conducted within the 28-day holding time for aqueous samples undergoing CLP protocol. Cyanide analyses were conducted within the 14-day holding time. The cooler temperature upon receipt at the laboratory was within the recommended temperature of 4°C +/- 2°C.

2. CALIBRATION

A. Initial Calibration

The percent recoveries for the Initial Calibration Verification (ICV) standard were within Quality Control (QC) limits for all constituents.

B. Continuing Calibration

The percent recoveries for the Continuing Calibration Verification (CCV) standard were within QC limits for all constituents.

3. BLANKS

The Initial Calibration Blank (ICB), Continuing Calibration Blanks (CCB) and Preparation Blanks (PB) were analyzed at the appropriate frequencies. No constituents were detected in the ICB, CCB, PB, Equipment blanks above the corresponding Contract Required Detection Limit (CRDL).

4. ICP INTERFERENCE CHECK SAMPLE

Results for the ICP analysis of the Interference Check Sample (ICS) solution AB were within 20% of the true value.

5. LABORATORY CONTROL SAMPLES

Recoveries were within the control limit (80-120%) for all constituents.

6. DUPLICATE ANALYSIS

The Relative Percent Difference (RPD) between the sample and duplicate results were within the acceptance criteria for all target compounds.

7. SPIKE SAMPLE ANALYSIS

The laboratory used sample SKSWD031009 for the matrix spike sample. The MS percent recoveries were within the acceptance criteria (75%-125%) with the exception of Selenium in the total (0%) and dissolved (50%) fractions. As per the National Functional Guidelines: if the percent recovery is greater than 30% and less than 74% qualify detected results for that analyte with "J" and non-detected results with "UJ". If the percent recovery is less than 30% qualify detected results for that analyte with "J" and non-detected results with "R".

8. ICP SERIAL DILUTION

As noted in the National Functional Guidelines: If the analyte concentration is at least 50 times above the IDL, its serial dilution analysis must then agree within 10% of the original determination after corrected for dilution. The serial dilution is performed to determine whether any significant chemical or physical interference's exist due to matrix effects. The serial dilution percent differences were within the acceptance criteria for all target analytes.

9. SYSTEM PERFORMANCE

The analytical system appears to have been working well at the time of these analyses, based on the evaluation of the raw data.

10. DOCUMENTATION

The documentation appeared accurate and in order.

11. OVERALL ASSESSMENT

The percent recoveries for Copper in the Contract Required Detection Limit (CRDL) standards were 63.0, 62.9%, and 69.1%. The detected Copper results greater than the IDL but less than two times the CRDL were qualified with as estimated with "J". The percent recoveries for Selenium in the Contract Required Detection Limit (CRDL) standards were 62.4 and 64.2%. The Selenium results were previously qualified under Section 7-titled "Spike Sample Analysis". The results are acceptable with the validator-added qualifiers.

**DATA VALIDATION SUMMARY – SAMPLE DELIVERY GROUP 204030804
SEMIVOLATILE ORGANICS**

Validation of the Gas Chromatograph/Mass Spectrometer (GC/MS) semi-volatile organics data, as prepared by Gulf Coast Analytical Laboratories (GCAL) for the samples collected from the Skinner Landfill site in March 2004, was conducted by Earth Tech using the National Functional Guidelines for Organic Data Review, (US EPA, October, 1999) as appropriate. The results were reported by GCAL under SDG 204030804.

| GCAL # | Sample Description |
|--------------|--------------------|
| 204030804-01 | SKSWD031009 |
| 204030804-02 | SKSWD03D1009 |
| 204030804-03 | SKSWD03MS1009 |
| 204030804-04 | SKSWD03MSD1009 |
| 204030804-06 | SKSWDEB1009 |

INTRODUCTION

Analyses were performed according to CLP-Organic Analysis Multi-Media, Multi-Concentration OLM04.2 SOW. Results of the sample analyses are reported by the laboratory as either qualified or unqualified. Unqualified results mean that the reported values may be used without reservation. Various data qualifier codes are used by the laboratory to denote specific information regarding the analytical results. The data validation process is intended to evaluate the data on a technical basis. The data package also was subjected to an internal laboratory quality review prior to submission to Earth Tech for data validation.

During the validation process, laboratory-qualified and unqualified data are verified against all available supporting documentation. Based on this evaluation, qualifier codes may be added, deleted or modified by the data user. Final results are therefore, either qualified or unqualified. Validator-qualified results are annotated with the following codes in accordance with the Functional Guidelines:

- U The constituent was analyzed for, but was not detected above the level of the associated analytical reporting limit. The associated value is either the sample quantitation limit or the sample detection limit.
- J The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample.
- UJ The analyte was not detected above the reported sample quantitation limit. However, the reported quantitation limit is approximate and may or may not represent the actual limit of quantitation necessary to accurately and precisely measure the analyte in the sample.

- R The sample results are rejected due to serious deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte cannot be verified.

Details of the semivolatile data validation findings and conclusions are provided in the following sections of this report:

1. Holding Times
2. GC/MS Tuning
3. Calibration
 - A. IC
 - B. CC
4. Blanks
5. System Monitoring Compound Recovery
6. MS/MSD
7. Internal Standards Performance
8. Compound Identification
9. Constituent Quantitation and Reported Detection Limits
10. System Performance
11. Documentation
12. Overall Assessment

1. HOLDING TIMES

All samples were initially extracted within the seven-day technical holding time and the five-day VTSR method holding time. The cooler temperature upon receipt at the laboratory was within the recommended temperature of $4^{\circ}\text{C} \pm 2^{\circ}\text{C}$.

2. GC/MS TUNING

The samples were analyzed on a single GC/MS system, identified as MSSV2. One decafluorotriphenylphosphine (DFTPP) tune was run representing the shift in which the standards and samples were analyzed. The DFTPP tune is acceptable.

3. CALIBRATION

A. Initial Calibration

One IC dated 3/19/04 was analyzed in support of the semivolatile sample analyses. Documentation of the IC was present in the data package, and the Relative Response Factor (RRF), as well as percent % RSD values were accurately reported for all target compounds. The criteria employed for technical data review purposes are different than those used in the method. The laboratory must meet a minimum RRF of 0.01; however, for data review purposes, a RRF criterion of "greater than or equal to 0.05" is applied to all semi-volatile compounds. The RRF's and the average RRF were within the acceptance criteria specified in the method for all reported analytes. The %RSD's were within the acceptance criteria specified in the method for all target analytes with the exception of Di-n-butylphthalate (32.0%), Di-n-octylphthalate (30.3%) and Diethylphthalate (41.8%). The lowest point of the calibration curve was dropped for Di-n-butylphthalate, Di-n-octylphthalate, and Diethylphthalate and the %RSD were recalculated. The recalculated %RSD were within the acceptance criteria of less than 30%. Di-n-butylphthalate, Di-n-octylphthalate, and Diethylphthalate were not detected in the associated samples therefore data qualification was not required.

B. Continuing Calibration

One CC dated 3/19/04 was analyzed in support of the semivolatile sample analyses reported in the data submissions. The RRF's for the CC dated 3/19/04 were within the acceptance criteria. The percent difference (%D) between the average RRF's and the CC Response Factors for the CC dated 3/19/04 were within the acceptance criteria with the exception the %D for Naphthalene. As per the National Functional Guidelines, if the %D exceeds the acceptance criteria qualify detected results for that analyte with "J" and non-detected results for that analyte with "UJ".

4. BLANKS

One laboratory semivolatile method blank and equipment blank were analyzed with this SDG. The results are summarized below.

Method Blank (0308SBLK)

There were not target analytes detected in method blank 0308SBLK.

Equipment Blank (SKSWDEB1009)

There were not target analytes detected in method blank SKSWDEB1009 collected on 3/2/04.

5. SYSTEM MONITORING COMPOUND RECOVERY

All reported semivolatile system monitoring compounds were recovered within acceptable control limits.

6. MATRIX SPIKE MATRIX SPIKE DUPLICATE (MS/MSD)

Sample SKSWD031009 was submitted for MS/MSD analysis. The MS/MSD percent recoveries were within the acceptance criteria with the exception of the 4-Nitrophenol. The %RPD between the MS/MSD were within the acceptance criteria. As per the National Functional Guidelines, no action is taken on MS/MSD data alone.

7. INTERNAL STANDARDS PERFORMANCE

Internal standard areas and retention times were within acceptable limits for the reported semivolatile sample analyses.

8. COMPOUND IDENTIFICATION

All reported semivolatile constituents were correctly identified with supporting chromatograms present in the data package.

9. CONSTITUENT QUANTITATION AND REPORTED DETECTION LIMITS

Constituent quantitations were correctly calculated and reported for semivolatile constituents.

10. SYSTEM PERFORMANCE

The analytical system appears to have been working well at the time of these analyses, based on the evaluation of the raw data submitted for review.

11. DOCUMENTATION

The documentation appeared accurate and in order with the exception of the following: GCAL inadvertently left the Date Extracted off of the CLP Form 1's therefore the data validator inserted the Date Extracted on the CLP Form 1's.

12. OVERALL ASSESSMENT

The results are acceptable with the validator-added qualifiers.

DATA VALIDATION SUMMARY – SAMPLE DELIVERY GROUP 204030804 VOLATILE ORGANIC

Validation of the GC/MS volatile organics data, as prepared by Gulf Coast Analytical Laboratories (GCAL) for the samples collected from the Skinner Landfill site in March 2004, was conducted by Earth Tech using the National Functional Guidelines for Organic Data Review, (US EPA, October, 1999), as appropriate. The results were reported by GCAL under SDG 204030804.

| GCAL # | Sample Description |
|--------------|--------------------|
| 204030804-01 | SKSWD031009 |
| 204030804-02 | SKSWD03D1009 |
| 204030804-03 | SKSWD03MS1009 |
| 204030804-04 | SKSWD03MSD1009 |
| 204030804-06 | SKSWDEB1009 |
| 204030804-07 | Trip Blank |

INTRODUCTION

Analyses were performed according to CLP-Organic Analysis Low Concentration OLC02.0 SOW. Results of the sample analyses are reported by the laboratory as either qualified or unqualified. Unqualified results mean that the reported values may be used without reservation. Various qualifier codes are used by the laboratory to denote specific information regarding the analytical results. The data validation process is intended to evaluate the data on a technical basis. The data package also was subjected to an internal laboratory quality review prior to submission to Earth Tech for data validation.

During the validation process, laboratory-qualified and unqualified data are verified against all available supporting documentation. Based on this evaluation, qualifier codes may be added, deleted or modified by the data user. Final results are therefore, either qualified or unqualified. Validator-qualified results are annotated with the following codes in accordance with the Functional Guidelines:

- U The constituent was analyzed for, but was not detected above the level of the associated analytical reporting limit. The associated value is either the sample quantitation limit or the sample detection limit.
- J The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample.
- UJ The analyte was not detected above the reported sample quantitation limit.

However, the reported quantitation limit is approximate and may or may not represent the actual limit of quantitation necessary to accurately and precisely measure the analyte in the sample.

- R** The sample results are rejected due to serious deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte cannot be verified.

The volatiles data validation findings and conclusions are provided in the following sections of this report:

1. Holding Times
2. GC/MS Tuning
3. Calibration
 - A. IC
 - B. CC
4. Blanks
5. System Monitoring Compound Recovery
6. MS/MSD
7. Laboratory Control Sample
8. Internal Standards Performance
9. Compound Identification
10. Constituent Quantitation and Reported Detection Limits
11. System Performance
12. Documentation
13. Overall Assessment

1. HOLDING TIMES

All samples for Volatile Organic Compounds (VOC) analyses were analyzed within the 14-day technical holding time and the 10-day VTSR method holding time. The cooler temperature upon receipt at the laboratory was within the recommended temperature of 4°C +/- 2°C.

2. GC/MS TUNING

All samples were analyzed on a single GC/MS system, identified as MSV2. One bromofluorobenzene (BFB) tune was run. The BFB tune is acceptable.

3. CALIBRATION

A. Initial Calibration

One IC dated 3/13/04 was analyzed on Instrument MSV2 in support of the volatile sample analyses reported in the data submissions. Documentation of the IC standards was present in the data package, and RRF's as well as %RSD values were accurately reported. The criteria employed for technical data review purposes are different than those used in the method. The laboratory must meet a minimum RRF of 0.01; however, for data review purposes, a RRF criterion of "greater than or equal to 0.05" is applied to all volatile compounds. The %RSD's were within the acceptance criteria specified in the method for all target analytes. The RRF's and the average RRF for the IC dated 3/13/04 were within the acceptance criteria specified in the method for all target analytes with the exception of Acetone. As per the National Functional Guidelines, if any initial calibration RRF is less than 0.05, qualify positive results that have acceptable mass spectral identification with "J", using professional judgement, and non-detected analytes as unusable (R).

B. Continuing Calibration

One CC dated 3/13/04 was analyzed on instrument MSV2 in support of the volatile sample analyses reported in the data submissions. The percent difference (%D) between the average RRF's and the CC RF's were within the acceptance criteria for all target analytes. The CC RRF's were within the acceptance criteria specified in the method for all target analytes with the exception of Acetone. The Acetone results were previously qualified under section 3A above.

4. BLANKS

One laboratory volatile method blank, storage blank, Trip Blank, and an Equipment Blank were analyzed with this SDG. The results are summarized below.

V2BLK01 (MB151783)

Methylene chloride, Chloroform and 1,3-Dichlorobenzene were detected at concentrations of 0.12 ppb, 0.38 ppb, and 0.034 ppb respectively in the method blank analyzed on 3/13/04.

Storage Blank (VHBLK01)

Methylene chloride (0.15 ppb) and 1,4-Dichlorobenzene (0.21 ppb) were detected in the storage blank analyzed on 3/13/04.

Trip Blank

Methylene chloride was detected at a concentration of 0.47 ppb in the Trip Blank submitted for the sampling event that occurred on 3/2/04. The Methylene chloride detected in the trip blank was mitigated by the presence of Methylene chloride in the associated method blank.

Equipment Blank (SKSWDEB1009)

Ethylbenzene (0.022 ppb), Methylene chloride (0.4 ppb), Styrene (0.06 ppb) Toluene (0.43 ppb) and total Xylenes (0.11 ppb) were detected in the Equipment Blank collected on 3/2/04. The analytes Ethylbenzene, Styrene, Toluene, and total Xylenes were not detected in the associated samples therefore no data qualification was not required. The Methylene chloride detected in the equipment blank was mitigated by the presence of Methylene chloride in the associated method blank.

5. SYSTEM MONITORING COMPOUND RECOVERY

All reported volatile system monitoring compounds were recovered within acceptable control limits for all samples.

6. MATRIX SPIKE/MATRIX SPIKE DUPLICATE

Sample SKSWD031009 was submitted for MS/MSD analysis. The MS/MSD percent recoveries and %RPD between the MS/MSD were within the acceptance criteria.

7. LABORATORY CONTROL SAMPLE

A LCS was analyzed in conjunction with this SDG. Recoveries were within the control limit for all constituents.

8. INTERNAL STANDARDS PERFORMANCE

Internal Standard areas and retention times were within acceptable limits for the reported volatile sample analyses.

9. COMPOUND IDENTIFICATION

All reported VOCs were correctly identified with supporting chromatograms present in the data package.

10. CONSTITUENT QUANTITATION AND REPORTED DETECTION LIMITS

Constituent quantitations were correctly calculated and reported for VOCs with the exception of Ethylbenzene. The Ethylbenzene standard and detected results were originally quantitated using the incorrect quantitation ion (GCAL used 106 instead of 91).

GCAL corrected the mistake and re-submitted the corrected pages that were affected in the laboratory report. The overall effect had no impact in the final result for Ethylbenzene.

11. SYSTEM PERFORMANCE

The analytical system appears to have been working well at the time of these analyses, based on the evaluation of the raw data.

12. DOCUMENTATION

The documentation appeared accurate and in order with the exception of the following: GCAL inadvertently left the "B" qualifier off of the CLP Form 1's for Methylene chloride therefore the data validator inserted a "B" qualifier in the "Q" column of the CLP Form 1's. The "B" qualifier indicates that the analyte was detected in the associated method blank.

13. OVERALL ASSESSMENT

The Acetone detected in sample SKSWD03D1009 could be do to low level contamination because Acetone is a common laboratory contaminant and the fact that Acetone was not detected in the associated duplicate sample. The results are acceptable with the validator-added qualifiers.

**DATA VALIDATION SUMMARY - SAMPLE DELIVERY GROUP 204030804
PESTICIDES**

Validation of the Gas Chromatography (GC) pesticides data, as prepared by Gulf Coast Analytical Laboratories (GCAL) for the samples collected from the Skinner Landfill site in March 2004, was conducted by Earth Tech using the National Functional Guidelines for Organic Data Review, (US EPA, October, 1999), as appropriate. The results were reported by GCAL under SDG 204030804.

| GCAL # | Sample Description |
|---------------|---------------------------|
| 204030804-01 | SKSWD031009 |
| 204030804-02 | SKSWD03D1009 |
| 204030804-03 | SKSWD03MS1009 |
| 204030804-04 | SKSWD03MSD1009 |
| 204030804-06 | SKSWDEB1009 |

INTRODUCTION

Analyses were performed according to CLP-Organic Analysis Multi-Media, Multi-Concentration OLM04.2 SOW. Results of the sample analyses are reported by the laboratory as either qualified or unqualified. Unqualified results mean that the reported values may be used without reservation. Various qualifier codes are used by the laboratory to denote specific information regarding the analytical results.

The data validation process is intended to evaluate the data on a technical basis. The data package also was subjected to an internal laboratory quality review prior to submission to Earth Tech for data validation.

During the validation process, laboratory-qualified and unqualified data are verified against all available supporting documentation. Based on this evaluation, qualifier codes may be added, deleted or modified by the data user. Final results are therefore, either qualified or unqualified. Validator-qualified results are annotated with the following codes in accordance with the Functional Guidelines:

- U** The constituent was analyzed for, but was not detected above the level of the associated analytical reporting limit. The associated value is either the sample quantitation limit or the sample detection limit.
- J** The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample.
- UJ** The analyte was not detected above the reported sample quantitation limit. However, the reported quantitation limit is approximate and may or may not represent the actual limit of quantitation necessary to accurately and precisely measure the analyte in the sample.

- R The sample results are rejected due to serious deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte cannot be verified.

Details of the pesticide data validation findings and conclusions are provided in the following sections of this report:

1. Holding Times
2. Gas Chromatograph/Electronic Capture Detector (GC/ECD) Instrument Performance Check
3. IC
4. Calibration Verification
5. Blanks
6. Surrogate Spikes
7. Matrix Spike/Matrix Spike Duplicate (MS/MSD)
8. Pesticide Cleanup Checks
9. Target Compound Identification
10. Constituent Quantitation and Reported Detection Limits
11. Documentation
12. Overall Assessment

1. HOLDING TIMES

All samples were extracted within the seven-day technical holding time and the five-day VTSR method holding time. The cooler temperature upon receipt at the laboratory was within the recommended temperature of 4°C +/- 2°C.

2. GC/ECD INSTRUMENT PERFORMANCE CHECK

The Performance Evaluation Mixture (PEM) was analyzed at the correct frequency. Absolute retention times were within limits. The percent resolution between adjacent peaks was within QC limits for the Pesticide Analyte Resolution Check. The percent resolution between adjacent peaks is within QC limits for the Performance Evaluation Mixtures (PEM).

The percent breakdown for both 4,4-DDT and Endrin in each PEM was less than 20.0% for both GC columns. The combined percent breakdown for 4,4-DDT and Endrin in each PEM was less than 30.0% for both GC columns.

3. INITIAL CALIBRATION

Individual standard mixtures A and B were analyzed at the correct frequencies and concentrations. The percent resolution criterion was met for Individual standard mixtures A and B.

The Percent Relative Standard Deviation (%RSD) of the calibration factors for each of the single component pesticides was less than 20%.

The multi-component target compounds were analyzed separately on both columns at a single concentration level. Retention times were determined from a minimum of three peaks.

4. CALIBRATION VERIFICATION

Absolute retention times were within appropriate time retention windows

5. BLANKS

One laboratory method blank and equipment blank were analyzed with this SDG. The results are summarized below.

Method Blank 151585

No constituents were detected above the laboratory-reporting limit. This blank corresponds to all samples extracted on 3/8/04.

Equipment Blank SKSWDEB1009

No constituents were detected above the laboratory-reporting limit in the equipment blank collected on 3/2/04.

6. SURROGATE SPIKES

Decachlorobiphenyl (DCB) and tetrachloro-m-xylene (TCX) surrogate spike recoveries were within the acceptance criteria for all samples except for TCX and DCB associated with sample SKSWD031009. There were no target compounds detected in sample SKSWD031009 therefore no action was taken.

7. MATRIX SPIKE/MATRIX SPIKE DUPLICATE

There were no samples submitted for MS MSD analysis during this sampling event.

8. PESTICIDE CLEANUP CHECKS

Recoveries of all pesticides and surrogates were within 80-120% for the lot of Florisil cartridges utilized for pesticide cleanup with the exception of Endrin (135%), 4,4'-DDT (139%) and Methoxychlor (121%). There were no target compounds detected in the associated samples therefore no action was taken.

9. TARGET COMPOUND IDENTIFICATION

All reported pesticide data were correctly identified with supporting chromatograms present in the data package.

10. CONSTITUENT QUANTITATION AND REPORTED DETECTION LIMITS

Constituent quantitations were correctly calculated and reported for pesticide constituents.

11. DOCUMENTATION

The documentation appeared accurate and in order.

12. OVERALL ASSESSMENT

The results are acceptable with the validator-added qualifiers.

REFERENCES

US EPA, 1994. *National Functional Guidelines for Inorganic Data Review.*

US EPA, 1999. *National Functional Guidelines for Organic Data Review.*



ANALYTICAL RESULTS

PERFORMED BY

GULF COAST ANALYTICAL LABORATORIES, INC.

Report Date 03/31/2004

GCAL Report 204030804

Deliver To Earth Tech
200 Vine Street
Wilder, KY 41076
859-442-2300

Attn Pat Higgins

Customer Earth Tech

Project Skinner Landfill

000001

CASE NARRATIVE

Client: Earth Tech **Report:** 204030804

Gulf Coast Analytical Laboratories received and analyzed the sample(s) listed on the sample cross-reference page of this report. Receipt of the sample(s) is documented by the attached chain of custody. This applies only to the sample(s) listed in this report. No sample integrity or quality control exceptions were identified unless noted below.

SEMI-VOLATILES MASS SPECTROMETRY

In the prep batch 270522, the MS/MSD exhibited sporadic recovery failures.

SEMI-VOLATILES GAS CHROMATOGRAPHY

In the Pesticide Florisil check analysis, the recoveries for Endrin, 4,4-DDT and Methoxychor were above recovery limits; however, these compounds were not detected in the associated samples.

In the Pesticide analysis for sample 20403080401 (SKSWD031009), the surrogate Tetrachloro-m-xylene and Decachlorobiphenyl are above the control limits.

METALS

In the ILM04.1 - CLP Metals analysis for prep batches 270882 and 270883, the MS recovery was outside the control limits for Selenium. The LCS recovery was within control limits. This indicates the analysis is in control and the sample is affected by matrix interference. A post-digestion spike was performed on the QC sample for these batches with a recovery of 88% and 99%.

000002

Laboratory Endorsement

Sample analysis was performed in accordance with approved methodologies provided by the Environmental Protection Agency or other recognized agencies. The samples and their corresponding extracts will be maintained for a period of 30 days unless otherwise arranged. Following this retention period the samples will be disposed in accordance with GCAL's Standard Operating Procedures.

Common Abbreviations Utilized in this Report

| | |
|--------------|--|
| ND | Indicates the result was Not Detected at the specified RDL |
| DO | Indicates the result was Diluted Out |
| MI | Indicates the result was subject to Matrix Interference |
| TNTC | Indicates the result was Too Numerous To Count |
| SUBC | Indicates the analysis was Sub-Contracted |
| FLD | Indicates the analysis was performed in the Field |
| PQL | Practical Quantitation Limit |
| MDL | Method Detection Limit |
| RDL | Reporting Detection Limit |
| 00:00 | Reported as a time equivalent to 12:00 AM |

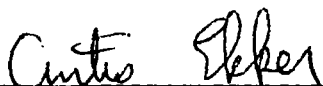
Reporting Flags Utilized in this Report

| | |
|----------|--|
| J | Indicates an estimated value |
| U | Indicates the compound was analyzed for but not detected |
| B | (ORGANICS) Indicates the analyte was detected in the associated Method Blank |
| B | (INORGANICS) Indicates the result is between the RDL and MDL |

Sample receipt at GCAL is documented through the attached chain of custody. In accordance with ISO Guide 25 and NELAC, this report shall be reproduced only in full and with the written permission of GCAL. The results contained within this report relate only to the samples reported. The documented results are presented within this report.

This report pertains only to the samples listed in the Report Sample Summary and should be retained as a permanent record thereof. The results contained within this report are intended for the use of the client. Any unauthorized use of the information contained in this report is prohibited.

I certify that this data package is in compliance with the terms and conditions of the contract and Statement of Work both technically and for completeness, for other than the conditions in the case narrative. Release of the data contained in this hardcopy data package and in the computer-readable data submitted has been authorized by the Quality Assurance Manager or his/her designee, as verified by the following signature.



CURTIS EKKER
DATA VALIDATION MANAGER
GCAL REPORT 204030804

THIS REPORT CONTAINS _____ PAGES.

000003

Report Sample Summary

| GCAL ID | Client ID | Matrix | Collect Date/Time | Receive Date/Time |
|-------------|----------------------|--------|-------------------|-------------------|
| 20403080401 | SKSWD031009 | Water | 03/02/2004 10:55 | 03/06/2004 11:50 |
| 20403080402 | SKSWD03D1009 | Water | 03/02/2004 11:18 | 03/06/2004 11:50 |
| 20403080403 | SKSWD03MS1009 | Water | 03/02/2004 11:45 | 03/06/2004 11:50 |
| 20403080404 | SKSWD03MSD1009 | Water | 03/02/2004 12:08 | 03/06/2004 11:50 |
| 20403080405 | SKSWD03DUP1009 | Water | 03/02/2004 12:08 | 03/06/2004 11:50 |
| 20403080406 | SKSWDEB1009 | Water | 03/02/2004 12:40 | 03/06/2004 11:50 |
| 20403080407 | TRIP BLANK | Water | 03/02/2004 00:00 | 03/06/2004 11:50 |
| 20403080408 | SKSWD031009 (DISS) | Water | 03/02/2004 10:55 | 03/06/2004 11:50 |
| 20403080409 | SKSWD03D1009 (DISS) | Water | 03/02/2004 11:18 | 03/06/2004 11:50 |
| 20403080410 | SKSWD03MS1009(DISS) | Water | 03/02/2004 11:45 | 03/06/2004 11:50 |
| 20403080411 | SKSWD03DUP1009(DISS) | Water | 03/02/2004 12:08 | 03/06/2004 11:50 |
| 20403080412 | SKSWDEB1009(DISS) | Water | 03/02/2004 12:40 | 03/06/2004 11:50 |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SKSWD031009

Lab Name: GCAL

Contract:

Lab Code: LA024

Case No.:

SAS No.:

SDG No.: 204030804

Matrix: (soil/water) Water

Sample wt/vol: 25 (g/ml) mL

Lab Sample ID: 20403080401

Level: (low/med)

Lab File ID: 2040313/T2457

% Moisture: not dec.

Date Collected: 03/02/04

Time: 1055

GC Column: DB-624-30M

ID: .53

(mm)

Date Received: 03/06/04

Instrument ID: MSV2

Date Analyzed: 03/13/04

Time: 1434

Soil Extract Volume: (µL)

Dilution Factor: 1

Analyst: RSP

Soil Aliquot Volume: (µL)

Prep Batch:

Analytical Batch: 270560

CONCENTRATION UNITS: ug/L

Analytical Method: OLC02.1 - CLP Vo

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|------------|---------------------------|----------|-----|-------|-----|
| 71-55-6 | 1,1,1-Trichloroethane | 1.0 | U | 0.010 | 1.0 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1.0 | U | 0.010 | 1.0 |
| 79-00-5 | 1,1,2-Trichloroethane | 1.0 | U | 0.010 | 1.0 |
| 75-34-3 | 1,1-Dichloroethane | 1.0 | U | 0.010 | 1.0 |
| 75-35-4 | 1,1-Dichloroethene | 1.0 | U | 0.010 | 1.0 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 106-93-4 | 1,2-Dibromoethane | 1.0 | U | 0.010 | 1.0 |
| 95-50-1 | 1,2-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 107-06-2 | 1,2-Dichloroethane | 1.0 | U | 0.010 | 1.0 |
| 540-59-0 | 1,2-Dichloroethene | 1.0 | U | 0.010 | 1.0 |
| 78-87-5 | 1,2-Dichloropropane | 1.0 | U | 0.010 | 1.0 |
| 541-73-1 | 1,3-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 106-46-7 | 1,4-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 78-93-3 | 2-Butanone | 5.0 | U | 0.010 | 5.0 |
| 591-78-6 | 2-Hexanone | 5.0 | U | 0.010 | 5.0 |
| 108-10-1 | 4-Methyl-2-pentanone | 5.0 | U | 0.010 | 5.0 |
| 67-64-1 | Acetone | 5.0 | U | 0.010 | 5.0 |
| 71-43-2 | Benzene | 1.0 | U | 0.010 | 1.0 |
| 75-27-4 | Bromodichloromethane | 1.0 | U | 0.010 | 1.0 |
| 75-25-2 | Bromoform | 1.0 | U | 0.010 | 1.0 |
| 74-83-9 | Bromomethane | 1.0 | U | 0.010 | 1.0 |
| 75-15-0 | Carbon disulfide | 1.0 | U | 0.010 | 1.0 |
| 56-23-5 | Carbon tetrachloride | 1.0 | U | 0.010 | 1.0 |
| 108-90-7 | Chlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 75-00-3 | Chloroethane | 1.0 | U | 0.010 | 1.0 |
| 67-66-3 | Chloroform | 1.0 | U | 0.010 | 1.0 |
| 74-87-3 | Chloromethane | 1.0 | U | 0.010 | 1.0 |
| 124-48-1 | Dibromochloromethane | 1.0 | U | 0.010 | 1.0 |
| 10061-01-5 | cis-1,3-Dichloropropene | 1.0 | U | 0.010 | 1.0 |
| 10061-02-6 | trans-1,3-Dichloropropene | 1.0 | U | 0.010 | 1.0 |
| 100-41-4 | Ethylbenzene | 1.0 | U | 0.010 | 1.0 |
| 75-09-2 | Methylene chloride | 2.0 0.12 | B/L | 0.010 | 2.0 |
| 100-42-5 | Styrene | 1.0 | U | 0.010 | 1.0 |

FORM 1 VOA

5/5/04
msk
000011

VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SKSWD031009

Lab Name: GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204030804
Matrix (soil/water): Water
Sample wt/vol: 25 (g/mL)
Level: (low/med) _____
% Moisture: not dec. _____
GC Column: DB-624-30M ID: .53 (mm)
Instrument ID: MSV2
Soil Extract Volume: _____ (µL)
Soil Aliquot Volume: _____ (µL)
Lab Sample ID: 20403080401
Lab File ID: 2040313/T2457
Date Collected: 03/02/04 Time: 1055
Date Received: 03/06/04
Date Analyzed: 03/13/04 Time: 1434
Dilution Factor: 1 Analyst: RSP
Prep Batch: _____ Analytical Batch: 270560
CONCENTRATION UNITS: µg/L
Analytical Method: OLC02.1 - CLP Vo

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|-----------|-------------------|--------|---|-------|-----|
| 127-18-4 | Tetrachloroethene | 1.0 | U | 0.010 | 1.0 |
| 108-88-3 | Toluene | 1.0 | U | 0.010 | 1.0 |
| 79-01-6 | Trichloroethene | 1.0 | U | 0.010 | 1.0 |
| 75-01-4 | Vinyl chloride | 1.0 | U | 0.010 | 1.0 |
| 1330-20-7 | Xylene (total) | 1.0 | U | 0.010 | 1.0 |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SKSWD03D1009

Lab Name: GCAL

Contract:

Lab Code: LA024

Case No.:

SAS No.:

SDG No.: 204030804

Matrix: (soil/water) Water

Sample wt/vol: 25 (g/ml) mL

Lab Sample ID: 20403080402

Level: (low/med)

Lab File ID: 2040313/T2456

% Moisture: not dec.

Date Collected: 03/02/04

Time: 1118

GC Column: DB-624-30M

ID: .53

(mm)

Date Received: 03/06/04

Instrument ID: MSV2

Date Analyzed: 03/13/04

Time: 1408

Soil Extract Volume: (µL)

Dilution Factor: 1

Analyst: RSP

Soil Aliquot Volume: (µL)

Prep Batch:

Analytical Batch: 270560

CONCENTRATION UNITS: ug/L

Analytical Method: OLC02.1 - CLP Vo

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|------------|---------------------------|----------|-------|-------|-----|
| 71-55-6 | 1,1,1-Trichloroethane | 1.0 | U | 0.010 | 1.0 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1.0 | U | 0.010 | 1.0 |
| 79-00-5 | 1,1,2-Trichloroethane | 1.0 | U | 0.010 | 1.0 |
| 75-34-3 | 1,1-Dichloroethane | 1.0 | U | 0.010 | 1.0 |
| 75-35-4 | 1,1-Dichloroethene | 1.0 | U | 0.010 | 1.0 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 106-93-4 | 1,2-Dibromoethane | 1.0 | U | 0.010 | 1.0 |
| 95-50-1 | 1,2-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 107-06-2 | 1,2-Dichloroethane | 1.0 | U | 0.010 | 1.0 |
| 540-59-0 | 1,2-Dichloroethene | 1.0 | U | 0.010 | 1.0 |
| 78-87-5 | 1,2-Dichloropropane | 1.0 | U | 0.010 | 1.0 |
| 541-73-1 | 1,3-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 106-46-7 | 1,4-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 78-93-3 | 2-Butanone | 5.0 | U | 0.010 | 5.0 |
| 591-78-6 | 2-Hexanone | 5.0 | U | 0.010 | 5.0 |
| 108-10-1 | 4-Methyl-2-pentanone | 5.0 | U | 0.010 | 5.0 |
| 67-64-1 | Acetone | 3.1 | J | 0.010 | 5.0 |
| 71-43-2 | Benzene | 1.0 | U | 0.010 | 1.0 |
| 75-27-4 | Bromodichloromethane | 1.0 | U | 0.010 | 1.0 |
| 75-25-2 | Bromoform | 1.0 | U | 0.010 | 1.0 |
| 74-83-9 | Bromomethane | 1.0 | U | 0.010 | 1.0 |
| 75-15-0 | Carbon disulfide | 1.0 | U | 0.010 | 1.0 |
| 56-23-5 | Carbon tetrachloride | 1.0 | U | 0.010 | 1.0 |
| 108-90-7 | Chlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 75-00-3 | Chloroethane | 1.0 | U | 0.010 | 1.0 |
| 67-66-3 | Chloroform | 1.0 | U | 0.010 | 1.0 |
| 74-87-3 | Chloromethane | 1.0 | U | 0.010 | 1.0 |
| 124-48-1 | Dibromochloromethane | 1.0 | U | 0.010 | 1.0 |
| 10061-01-5 | cis-1,3-Dichloropropene | 1.0 | U | 0.010 | 1.0 |
| 10061-02-6 | trans-1,3-Dichloropropene | 1.0 | U | 0.010 | 1.0 |
| 100-41-4 | Ethylbenzene | 1.0 | U | 0.010 | 1.0 |
| 75-09-2 | Methylene chloride | 2.0 2.28 | 3.4 4 | 0.010 | 2.0 |
| 100-42-5 | Styrene | 1.0 | U | 0.010 | 1.0 |

FORM 1 VOA

000017

5/5/04
msc

VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SKSWD03D1009

Lab Name: GCAL

Contract: _____

Lab Code: LA024

Case No.: _____

SAS No.: _____

SOG No.: 204030804

Matrix (soil/water): Water

Sample wt/vol: 25 (g/mL) mL

Lab Sample ID: 20403080402

Level: (low/med) _____

Lab File ID: 2040313/T2456

% Moisture: not dec. _____

Date Collected: 03/02/04

Time: 1118

GC Column: DB-624-30M ID: .53 (mm)

Date Received: 03/06/04

Instrument ID: MSV2

Date Analyzed: 03/13/04

Time: 1408

Soil Extract Volume: _____ (µL)

Dilution Factor: 1

Analyst: RSP

Soil Aliquot Volume: _____ (µL)

Prep Batch: _____

Analytical Batch: 270560

CONCENTRATION UNITS: ug/L

Analytical Method: OLC02.1 - CLP Vo

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|-----------|-------------------|--------|---|-------|-----|
| 127-18-4 | Tetrachloroethene | ND | U | 0.010 | 1.0 |
| 108-88-3 | Toluene | ND | U | 0.010 | 1.0 |
| 79-01-6 | Trichloroethene | ND | U | 0.010 | 1.0 |
| 75-01-4 | Vinyl chloride | ND | U | 0.010 | 1.0 |
| 1330-20-7 | Xylene (total) | ND | U | 0.010 | 1.0 |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SKSWDEB1009

Lab Name: GCAL Contract: _____
 Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204030804
 Matrix: (soil/water) Water
 Sample wt/vol: 25 (g/ml) mL Lab Sample ID: 20403080406
 Level: (low/med) _____ Lab File ID: 2040313/T2458
 % Moisture: not dec. _____ Date Collected: 03/02/04 Time: 1240
 GC Column: DB-624-30M ID: .53 (mm) Date Received: 03/06/04
 Instrument ID: MSV2 Date Analyzed: 03/13/04 Time: 1459
 Soil Extract Volume: _____ (μL) Dilution Factor: 1 Analyst: RSP
 Soil Aliquot Volume: _____ (μL) Prep Batch: _____ Analytical Batch: 270560
 Analytical Method: OLC02.1 - CLP Vo

CONCENTRATION UNITS: ug/L

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|------------|---------------------------|--------|---|-------|-----|
| 71-55-6 | 1,1,1-Trichloroethane | 1.0 | U | 0.010 | 1.0 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1.0 | U | 0.010 | 1.0 |
| 79-00-5 | 1,1,2-Trichloroethane | 1.0 | U | 0.010 | 1.0 |
| 75-34-3 | 1,1-Dichloroethane | 1.0 | U | 0.010 | 1.0 |
| 75-35-4 | 1,1-Dichloroethene | 1.0 | U | 0.010 | 1.0 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 106-93-4 | 1,2-Dibromoethane | 1.0 | U | 0.010 | 1.0 |
| 95-50-1 | 1,2-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 107-06-2 | 1,2-Dichloroethane | 1.0 | U | 0.010 | 1.0 |
| 540-59-0 | 1,2-Dichloroethene | 1.0 | U | 0.010 | 1.0 |
| 78-87-5 | 1,2-Dichloropropane | 1.0 | U | 0.010 | 1.0 |
| 541-73-1 | 1,3-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 106-46-7 | 1,4-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 78-93-3 | 2-Butanone | 5.0 | U | 0.010 | 5.0 |
| 591-78-6 | 2-Hexanone | 5.0 | U | 0.010 | 5.0 |
| 108-10-1 | 4-Methyl-2-pentanone | 5.0 | U | 0.010 | 5.0 |
| 67-64-1 | Acetone | 5.0 | U | 0.010 | 5.0 |
| 71-43-2 | Benzene | 1.0 | U | 0.010 | 1.0 |
| 75-27-4 | Bromodichloromethane | 1.0 | U | 0.010 | 1.0 |
| 75-25-2 | Bromoform | 1.0 | U | 0.010 | 1.0 |
| 74-83-9 | Bromomethane | 1.0 | U | 0.010 | 1.0 |
| 75-15-0 | Carbon disulfide | 1.0 | U | 0.010 | 1.0 |
| 56-23-5 | Carbon tetrachloride | 1.0 | U | 0.010 | 1.0 |
| 108-90-7 | Chlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 75-00-3 | Chloroethane | 1.0 | U | 0.010 | 1.0 |
| 67-66-3 | Chloroform | 1.0 | U | 0.010 | 1.0 |
| 74-87-3 | Chloromethane | 1.0 | U | 0.010 | 1.0 |
| 124-48-1 | Dibromochloromethane | 1.0 | U | 0.010 | 1.0 |
| 10061-01-5 | cis-1,3-Dichloropropene | 1.0 | U | 0.010 | 1.0 |
| 10061-02-6 | trans-1,3-Dichloropropene | 1.0 | U | 0.010 | 1.0 |
| 100-41-4 | Ethylbenzene | 0.022 | J | 0.010 | 1.0 |
| 75-09-2 | Methylene chloride | 2.0240 | B | 0.010 | 2.0 |
| 100-42-5 | Styrene | 0.060 | J | 0.010 | 1.0 |

FORM 1 VOA

000024

5/16/04
msu

VOLATILE ORGANIC ANALYSIS DATA SHEET

SAMPLE NO.

SKSWDEB1009

Lab Name: GCAL

Contract

Lab Code: LA024

Case No:

SAS No:

SDG No: 204030804

Matrix (soil/water) Water

Sample wt/vol: 25 (g/ml) mL

Lab Sample ID: 20403080406

Level: (low/med)

Lab File ID: 2040313/T2458

% Moisture: not dec.

Date Collected: 03/02/04

Time: 1240

GC Column: DB-624-30M

ID: 53

(mm)

Date Received: 03/06/04

Instrument ID: MSV2

Date Analyzed: 03/13/04

Time: 1459

Soil Extract Volume: (µL)

Dilution Factor: 1

Analyst: RSP

Soil Aliquot Volume: (µL)

Prep Batch:

Analytical Batch: 270560

CONCENTRATION UNITS: µg/L

Analytical Method: OLC02.1 - CLP Vo

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|-----------|-------------------|--------|---|-------|-----|
| 127-18-4 | Tetrachloroethene | 1.0 | U | 0.010 | 1.0 |
| 106-68-3 | Toluene | 0.43 | J | 0.010 | 1.0 |
| 79-01-6 | Trichloroethene | 1.0 | U | 0.010 | 1.0 |
| 75-01-4 | Vinyl chloride | 1.0 | U | 0.010 | 1.0 |
| 1330-20-7 | Xylene (total) | 0.11 | J | 0.010 | 1.0 |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

TRIP BLANK

Lab Name: GCAL Contract: _____

Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204030804

Matrix: (soil/water) Water

Sample wt/vol: 25 (g/ml) mL Lab Sample ID: 20403080407

Level: (low/med) _____ Lab File ID: 2040313/T2459

% Moisture: not dec. _____ Date Collected: 03/02/04 Time: 0000

GC Column: DB-624-30M ID: .53 (mm) Date Received: 03/06/04

Instrument ID: MSV2 Date Analyzed: 03/13/04 Time: 1524

Soil Extract Volume: _____ (μL) Dilution Factor: 1 Analyst: RSP

Soil Aliquot Volume: _____ (μL) Prep Batch: _____ Analytical Batch: 270560

CONCENTRATION UNITS: ug/L Analytical Method: OLC02.1 - CLP Vo

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|------------|---------------------------|----------|----|-------|-----|
| 71-55-6 | 1,1,1-Trichloroethane | 1.0 | U | 0.010 | 1.0 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1.0 | U | 0.010 | 1.0 |
| 79-00-5 | 1,1,2-Trichloroethane | 1.0 | U | 0.010 | 1.0 |
| 75-34-3 | 1,1-Dichloroethane | 1.0 | U | 0.010 | 1.0 |
| 75-35-4 | 1,1-Dichloroethene | 1.0 | U | 0.010 | 1.0 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 106-93-4 | 1,2-Dibromoethane | 1.0 | U | 0.010 | 1.0 |
| 95-50-1 | 1,2-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 107-06-2 | 1,2-Dichloroethane | 1.0 | U | 0.010 | 1.0 |
| 540-59-0 | 1,2-Dichloroethene | 1.0 | U | 0.010 | 1.0 |
| 78-87-5 | 1,2-Dichloropropane | 1.0 | U | 0.010 | 1.0 |
| 541-73-1 | 1,3-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 106-46-7 | 1,4-Dichlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 78-93-3 | 2-Butanone | 5.0 | U | 0.010 | 5.0 |
| 591-78-6 | 2-Hexanone | 5.0 | U | 0.010 | 5.0 |
| 108-10-1 | 4-Methyl-2-pentanone | 5.0 | U | 0.010 | 5.0 |
| 67-64-1 | Acetone | 5.0 | U | 0.010 | 5.0 |
| 71-43-2 | Benzene | 1.0 | U | 0.010 | 1.0 |
| 75-27-4 | Bromodichloromethane | 1.0 | U | 0.010 | 1.0 |
| 75-25-2 | Bromoform | 1.0 | U | 0.010 | 1.0 |
| 74-83-9 | Bromomethane | 1.0 | U | 0.010 | 1.0 |
| 75-15-0 | Carbon disulfide | 1.0 | U | 0.010 | 1.0 |
| 56-23-5 | Carbon tetrachloride | 1.0 | U | 0.010 | 1.0 |
| 108-90-7 | Chlorobenzene | 1.0 | U | 0.010 | 1.0 |
| 75-00-3 | Chloroethane | 1.0 | U | 0.010 | 1.0 |
| 67-66-3 | Chloroform | 1.0 | U | 0.010 | 1.0 |
| 74-87-3 | Chloromethane | 1.0 | U | 0.010 | 1.0 |
| 124-48-1 | Dibromochloromethane | 1.0 | U | 0.010 | 1.0 |
| 10061-01-5 | cis-1,3-Dichloropropene | 1.0 | U | 0.010 | 1.0 |
| 10061-02-6 | trans-1,3-Dichloropropene | 1.0 | U | 0.010 | 1.0 |
| 100-41-4 | Ethylbenzene | 1.0 | U | 0.010 | 1.0 |
| 75-09-2 | Methylene chloride | 2.0 0.47 | BS | 0.010 | 2.0 |
| 100-42-5 | Styrene | 1.0 | U | 0.010 | 1.0 |

VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

TRIP BLANK

Lab Name: GCAL

Contract: _____

Lab Code: LA024

Case No.: _____

SAS No.: _____

SDG No.: 204030804

Matrix (soil/water) Water

Sample wt/vol: 25 (g/ml) mL

Lab Sample ID: 20403080407

Level: (low/med) _____

Lab File ID: 2040313/T2459

% Moisture: not dec. _____

Date Collected: 03/02/04

Time: 0000

GC Column: DB-624-30M

ID: .53

(m)

Date Received: 03/06/04

Instrument ID: MSV2

Date Analyzed: 03/13/04

Time: 1524

Soil Extract Volume: _____ (µL)

Dilution Factor: 1

Analyst: RSP

Soil Aliquot Volume: _____ (µL)

Prep Batch: _____

Analytical Batch: 270560

CONCENTRATION UNITS µg/L

Analytical Method: OLC02.1 - CLP Vo

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|-----------|-------------------|--------|---|-------|-----|
| 127-18-4 | Tetrachloroethene | 1.0 | U | 0.010 | 1.0 |
| 108-88-3 | Toluene | 1.0 | U | 0.010 | 1.0 |
| 79-01-6 | Trichloroethene | 1.0 | U | 0.010 | 1.0 |
| 75-01-4 | Vinyl chloride | 1.0 | U | 0.010 | 1.0 |
| 1330-20-7 | Xylene (total) | 1.0 | U | 0.010 | 1.0 |

SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

Lab Name: GCAL Sample ID: SKSWD031909
 Lab Code: LA024 Case No.: _____ Contract: _____
 SAS No.: _____ SDG No.: 204030804 Lab File ID: 2040319/S4997
 Matrix: Water Lab Sample ID: 20403080401
 Sample wt/vol: 1000 Units: mL Date Collected: 03/02/04 Time: 1055
 Level: (low/med) _____ Date Received: 03/06/04
 % Moisture: _____ decanted: (Y/N) _____ Date Extracted: 3/8/04
 GC Column: DB-5MS-30M ID: .25 (mm) Date Analyzed: 03/19/04 Time: 1323
 Concentrated Extract Volume: 1000 (µL) Dilution Factor: 1 Analyst: RLW
 Injection Volume: 1.0 (µL) Prep Method: _____
 GPC Cleanup: (Y/N) N pH: _____ Analytical Method: OLMO 4.2
 Instrument ID: MSSV2

CONCENTRATION UNITS: ug/LPrep Batch: 270522 Analytical Batch: 271099

| CAS NO. | COMPOUND | RESULT | Q | MDL | PQL |
|-----------|-----------------------------|--------|---|-------|------|
| 95-95-4 | 2,4,5-Trichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 88-06-2 | 2,4,6-Trichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 120-83-2 | 2,4-Dichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 51-28-5 | 2,4-Dinitrophenol | 25.0 | U | 0.010 | 25.0 |
| 121-14-2 | 2,4-Dinitrotoluene | 10.0 | U | 0.010 | 10.0 |
| 606-20-2 | 2,6-Dinitrotoluene | 10.0 | U | 0.010 | 10.0 |
| 91-58-7 | 2-Chloronaphthalene | 10.0 | U | 0.010 | 10.0 |
| 95-57-8 | 2-Chlorophenol | 10.0 | U | 0.010 | 10.0 |
| 91-57-6 | 2-Methylnaphthalene | 10.0 | U | 0.010 | 10.0 |
| 88-74-4 | 2-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 88-75-5 | 2-Nitrophenol | 10.0 | U | 0.010 | 10.0 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 10.0 | U | 0.010 | 10.0 |
| 99-09-2 | 3-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 534-52-1 | 2-Methyl-4,6-dinitrophenol | 25.0 | U | 0.010 | 25.0 |
| 59-50-7 | 4-Chloro-3-methylphenol | 10.0 | U | 0.010 | 10.0 |
| 106-47-8 | 4-Chloroaniline | 10.0 | U | 0.010 | 10.0 |
| 7005-72-3 | 4-Chlorophenyl-phenylether | 10.0 | U | 0.010 | 10.0 |
| 106-44-5 | 4-Methylphenol (p-Cresol) | 10.0 | U | 0.010 | 10.0 |
| 83-32-9 | Acenaphthene | 10.0 | U | 0.010 | 10.0 |
| 208-96-8 | Acenaphthylene | 10.0 | U | 0.010 | 10.0 |
| 120-12-7 | Anthracene | 10.0 | U | 0.010 | 10.0 |
| 56-55-3 | Benzo(a)anthracene | 10.0 | U | 0.010 | 10.0 |
| 50-32-8 | Benzo(a)pyrene | 10.0 | U | 0.010 | 10.0 |
| 205-99-2 | Benzo(b)fluoranthene | 10.0 | U | 0.010 | 10.0 |
| 191-24-2 | Benzo(g,h,i)perylene | 10.0 | U | 0.010 | 10.0 |
| 207-08-9 | Benzo(k)fluoranthene | 10.0 | U | 0.010 | 10.0 |
| 111-91-1 | Bis(2-Chloroethoxy)methane | 10.0 | U | 0.010 | 10.0 |
| 111-44-4 | Bis(2-Chloroethyl)ether | 10.0 | U | 0.010 | 10.0 |
| 108-60-1 | bis(2-Chloroisopropyl)ether | 10.0 | U | 0.010 | 10.0 |
| 117-81-7 | bis(2-ethylhexyl)phthalate | 10.0 | U | 0.010 | 10.0 |

SEMI-VOLATILE ORGANICS ANALYSIS DATA SHEET

Lab Name: GCAL
 Lab Code: LA024 Case No.: _____
 SAS No.: _____ SDG No.: 204030804
 Matrix: Water
 Sample wt/vol: 1000 Units: µL
 Level: (lowmed) _____
 % Moisture: _____ decanted: (Y/N) _____
 GC Column: DB-SMS-30M ID: 25 (mm)
 Concentrated Extract Volume: 1000 (µL)
 Injection Volume: 1.0 (µL)
 GPC Cleanup: (Y/N) N pH: _____

Sample ID: SKS:WD031009
 Contract: _____
 Lab File ID: 2040319/S4997
 Lab Sample ID: 20403080401
 Date Collected: 03/02/04 Time: 1055
 Date Received: 03/06/04
 Date Extracted: 3/5/04
 Date Analyzed: 03/19/04 Time: 1323
 Dilution Factor: 1 Analyst: RLW
 Prep Method: _____
 Analytical Method: OLMO 4.2
 Instrument ID: MSSV2

CONCENTRATION UNITS: µg/L

Prep Batch: 270522 Analytical Batch: 271099

| CAS NO. | COMPOUND | RESULT | Q | MDL | PQL |
|----------|---------------------------|--------|---|-------|------|
| 101-55-3 | 4-Bromophenyl-phenylether | 10.0 | U | 0.010 | 10.0 |
| 85-68-7 | Butylbenzylphthalate | 10.0 | U | 0.010 | 10.0 |
| 86-74-8 | Carbazole | 10.0 | U | 0.010 | 10.0 |
| 218-01-9 | Chrysene | 10.0 | U | 0.010 | 10.0 |
| 84-74-2 | Di-n-butylphthalate | 10.0 | U | 0.010 | 10.0 |
| 117-84-0 | Di-n-octylphthalate | 10.0 | U | 0.010 | 10.0 |
| 53-70-3 | Dibenz[a,h]anthracene | 10.0 | U | 0.010 | 10.0 |
| 132-64-9 | Dibenzofuran | 10.0 | U | 0.010 | 10.0 |
| 84-86-2 | Diethylphthalate | 10.0 | U | 0.010 | 10.0 |
| 131-11-3 | Dimethyl-phthalate | 10.0 | U | 0.010 | 10.0 |
| 105-67-9 | 2,4-Dimethylchloro | 10.0 | U | 0.010 | 10.0 |
| 206-44-0 | Fluoranthene | 10.0 | U | 0.010 | 10.0 |
| 96-73-7 | Fluorene | 10.0 | U | 0.010 | 10.0 |
| 118-74-1 | Hexachlorobenzene | 10.0 | U | 0.010 | 10.0 |
| 87-68-3 | Hexachlorobutadiene | 10.0 | U | 0.010 | 10.0 |
| 77-47-4 | Hexachlorocyclopentadiene | 10.0 | U | 0.010 | 10.0 |
| 67-72-1 | Hexachloroethane | 10.0 | U | 0.010 | 10.0 |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | 10.0 | U | 0.010 | 10.0 |
| 78-59-1 | Isophorone | 10.0 | U | 0.010 | 10.0 |
| 91-20-3 | Naphthalene | 10.0 | U | 0.010 | 10.0 |
| 100-01-6 | 4-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 98-95-3 | Nitrobenzene | 10.0 | U | 0.010 | 10.0 |
| 100-02-7 | 4-Nitrophenol | 25.0 | U | 0.010 | 25.0 |
| 87-86-5 | Pentachlorobenzol | 25.0 | U | 0.010 | 25.0 |
| 85-01-8 | Phenanthrene | 10.0 | U | 0.010 | 10.0 |
| 108-95-2 | Phenol | 10.0 | U | 0.010 | 10.0 |
| 129-00-0 | Pyrene | 10.0 | U | 0.010 | 10.0 |
| 621-64-7 | N-Nitroso-dim-propylamine | 10.0 | U | 0.010 | 10.0 |
| 86-30-6 | N-Nitrosodiphenylamine | 10.0 | U | 0.010 | 10.0 |
| 95-48-7 | o-Cresol | 10.0 | U | 0.010 | 10.0 |

1F
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

Lab Name: GCAL Sample ID: SKSWD031009
Lab Code: LA024 2 Case No.: _____ Contract: _____
SAS No.: _____ SDG No.: 204030804 Lab File ID: 2040319/S4997
Matrix: Water Lab Sample ID: 204030804C1
Sample wt/vol: _____ Units: _____ Date Collected: 03/02/04 Time: 1055
Level: (low/med) _____ Date Received: 03/06/04
% Moisture: not dec. _____ Date Extracted: 3/8/04
GC Column: DB-5MS-30M ID: .25 (mm) Date Analyzed: 03/19/04 Time: 1323
Concentrated Extract Volume: 1000 (µL) Dilution Factor: 1 Analyst: RLW
Injection Volume: 1.0 (µL) Prep Method: _____
GPC Cleanup: (Y/N) N pH: _____ Analytical Method: SW-646 8270C
Instrument ID: MSSV2

Number TICs Found : 3

CONCENTRATION UNITS:

| | CAS NO. | COMPOUND | RT | EST. CONC. | Q |
|----|-----------|--------------------------------|-------|------------|---|
| 1. | | Unknown | 4.364 | 3.38 | |
| 2. | | Unknown | 4.625 | 10.5 | |
| 3. | 4291-79-6 | Cyclohexane, 1-methyl-2-propyl | 4.862 | 13.5 | |

5/16/04
RLW

SEMIVOLATILE ORGANIC ANALYSIS DATA SHEET

Lab Name: GCAL Sample ID: SKSWD03D1009
 Lab Code: LA024 Case No.: _____ Contract: _____
 SAS No.: _____ SOG No.: 204030804 Lab File ID: 2040319/S4998
 Matrix: Water Lab Sample ID: 20403080402
 Sample wt/vol: 1000 Units mL Date Collected: 03/02/04 Time: 1118
 Level: (low/med) _____ Date Received: 03/06/04
 % Moisture: _____ deaerated (Y/N) _____ Date Extracted: 3/16/04
 GC Column: DB-5MS-30M ID: 25 (mm) Date Analyzed: 03/19/04 Time: 1351
 Concentrated Extract Volume: 1000 (µL) Dilution Factor: 1 Analyst: RLW
 Injection Volume: 1.0 (µL) Prep Method: _____
 GPC Cleanup: (Y/N) N pH: _____ Analytical Method: OLMO 4.2

CONCENTRATION UNITS: µg/L

Instrument ID: MSSV2
 Prep Batch: 270522 Analytical Batch: 271099

| CAS NO. | COMPOUND | RESULT | Q | MDL | PQL |
|-----------|-----------------------------|--------|---|-------|------|
| 95-95-4 | 2,4,5-Trichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 88-06-2 | 2,4,6-Trichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 120-83-2 | 2,4-Dichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 51-28-5 | 2,4-Dinitrophenol | 25.0 | U | 0.010 | 25.0 |
| 121-14-2 | 2,4-Dinitrotoluene | 10.0 | U | 0.010 | 10.0 |
| 606-20-2 | 2,6-Dinitrotoluene | 10.0 | U | 0.010 | 10.0 |
| 91-58-7 | 2-Chloronaphthalene | 10.0 | U | 0.010 | 10.0 |
| 95-57-8 | 2-Chlorophenol | 10.0 | U | 0.010 | 10.0 |
| 91-57-6 | 2-Methylnaphthalene | 10.0 | U | 0.010 | 10.0 |
| 88-74-4 | 2-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 88-75-5 | 2-Nitrophenol | 10.0 | U | 0.010 | 10.0 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 10.0 | U | 0.010 | 10.0 |
| 99-08-2 | 3-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 534-52-1 | 2-Methyl-4,6-dinitrophenol | 25.0 | U | 0.010 | 25.0 |
| 59-50-7 | 4-Chloro-3-methylphenol | 10.0 | U | 0.010 | 10.0 |
| 106-47-8 | 4-Chloroaniline | 10.0 | U | 0.010 | 10.0 |
| 7005-72-3 | 4-Chlorophenylphenylether | 10.0 | U | 0.010 | 10.0 |
| 106-44-5 | 4-Methylphenol (p-Cresol) | 10.0 | U | 0.010 | 10.0 |
| 83-32-9 | Acenaphthene | 10.0 | U | 0.010 | 10.0 |
| 208-96-8 | Acenaphthylene | 10.0 | U | 0.010 | 10.0 |
| 120-12-7 | Anthracene | 10.0 | U | 0.010 | 10.0 |
| 56-55-3 | Benzo(a)anthracene | 10.0 | U | 0.010 | 10.0 |
| 50-32-8 | Benzo(a)pyrene | 10.0 | U | 0.010 | 10.0 |
| 205-99-2 | Benzo(b)fluoranthene | 10.0 | U | 0.010 | 10.0 |
| 191-24-2 | Benzo(g,h,i)perylene | 10.0 | U | 0.010 | 10.0 |
| 207-08-9 | Benzo(k)fluoranthene | 10.0 | U | 0.010 | 10.0 |
| 111-91-1 | Bis(2-Chloroethoxy)methane | 10.0 | U | 0.010 | 10.0 |
| 111-44-4 | Bis(2-Chloroethyl)ether | 10.0 | U | 0.010 | 10.0 |
| 108-60-1 | bis(2-Chloroisopropyl)ether | 10.0 | U | 0.010 | 10.0 |
| 117-81-7 | bis(2-ethylhexyloxy)ethane | 10.0 | U | 0.010 | 10.0 |

SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

Lab Name: GCAL Sample ID: SKSWD03D1009
 Lab Code: LA024 Case No.: _____ Contract: _____
 SAS No.: _____ SDG No.: 204030804 Lab File ID: 2040319/S4998
 Matrix: Water Lab Sample ID: 20403080402
 Sample wt/vol: 1000 Units: mL Date Collected: 03/02/04 Time: 1118
 Level: (low/med) _____ Date Received: 03/06/04
 % Moisture: _____ decanted: (Y/N) _____ Date Extracted: 3/8/04
 GC Column: DB-5MS-30M ID: .25 (mm) Date Analyzed: 03/19/04 Time: 1351
 Concentrated Extract Volume: 1000 (µL) Dilution Factor: 1 Analyst: RLW
 Injection Volume: 1.0 (µL) Prep Method: _____
 GPC Cleanup: (Y/N) N pH: _____ Analytical Method: OLMO 4.2
 Instrument ID: MSSV2
 Prep Batch: 270522 Analytical Batch: 271099

CONCENTRATION UNITS: ug/L

| CAS NO. | COMPOUND | RESULT | Q | MDL | PQL |
|----------|----------------------------|--------|---|-------|------|
| 101-55-3 | 4-Bromophenyl-phenylether | 10.0 | U | 0.010 | 10.0 |
| 85-68-7 | Butylbenzylphthalate | 10.0 | U | 0.010 | 10.0 |
| 86-74-8 | Carbazole | 10.0 | U | 0.010 | 10.0 |
| 218-01-9 | Chrysene | 10.0 | U | 0.010 | 10.0 |
| 84-74-2 | Di-n-butylphthalate | 10.0 | U | 0.010 | 10.0 |
| 117-84-0 | Di-n-octylphthalate | 10.0 | U | 0.010 | 10.0 |
| 53-70-3 | Dibenz(a,h)anthracene | 10.0 | U | 0.010 | 10.0 |
| 132-64-9 | Dibenzofuran | 10.0 | U | 0.010 | 10.0 |
| 84-66-2 | Diethylphthalate | 10.0 | U | 0.010 | 10.0 |
| 131-11-3 | Dimethyl-phthalate | 10.0 | U | 0.010 | 10.0 |
| 105-67-9 | 2,4-Dimethylphenol | 10.0 | U | 0.010 | 10.0 |
| 206-44-0 | Fluoranthene | 10.0 | U | 0.010 | 10.0 |
| 86-73-7 | Fluorene | 10.0 | U | 0.010 | 10.0 |
| 118-74-1 | Hexachlorobenzene | 10.0 | U | 0.010 | 10.0 |
| 87-68-3 | Hexachlorobutadiene | 10.0 | U | 0.010 | 10.0 |
| 77-47-4 | Hexachlorocyclopentadiene | 10.0 | U | 0.010 | 10.0 |
| 67-72-1 | Hexachloroethane | 10.0 | U | 0.010 | 10.0 |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | 10.0 | U | 0.010 | 10.0 |
| 78-59-1 | Isophorone | 10.0 | U | 0.010 | 10.0 |
| 91-20-3 | Naphthalene | 10.0 | U | 0.010 | 10.0 |
| 100-01-6 | 4-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 98-95-3 | Nitrobenzene | 10.0 | U | 0.010 | 10.0 |
| 100-02-7 | 4-Nitrophenol | 25.0 | U | 0.010 | 25.0 |
| 87-86-5 | Pentachlorophenol | 25.0 | U | 0.010 | 25.0 |
| 85-01-8 | Phenanthrene | 10.0 | U | 0.010 | 10.0 |
| 108-95-2 | Phenol | 10.0 | U | 0.010 | 10.0 |
| 129-00-0 | Pyrene | 10.0 | U | 0.010 | 10.0 |
| 621-64-7 | N-Nitroso-di-n-propylamine | 10.0 | U | 0.010 | 10.0 |
| 86-30-6 | N-Nitrosodiphenylamine | 10.0 | U | 0.010 | 10.0 |
| 95-48-7 | o-Cresol | 10.0 | U | 0.010 | 10.0 |

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SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

| | |
|--|---|
| Lab Name: <u>GCAL</u> | Sample ID: <u>SKSWD03D1009</u> |
| Lab Code: <u>LA024</u> 2 Case No.: _____ | Contract: _____ |
| SAS No.: _____ SDG No.: <u>204030804</u> | Lab File ID: <u>2040319/S4998</u> |
| Matrix: <u>Water</u> | Lab Sample ID: <u>20403080402</u> |
| Sample wt/vol: _____ Units: _____ | Date Collected: <u>03/02/04</u> Time: <u>1118</u> |
| Level: (low/med) _____ | Date Received: <u>03/06/04</u> |
| % Moisture: not dec. _____ | Date Extracted: <u>3/8/04</u> |
| GC Column: <u>DB-5MS-30M</u> ID: <u>.25</u> (mm) | Date Analyzed: <u>03/19/04</u> Time: <u>1351</u> |
| Concentrated Extract Volume: <u>1000</u> (µL) | Dilution Factor: <u>1</u> Analyst: <u>RLW</u> |
| Injection Volume: _____ (µL) | Prep Method: _____ |
| GPC Cleanup: (Y/N) <u>N</u> pH: _____ | Analytical Method: <u>SW-846 8270C</u> |
| | Instrument ID: <u>MSSV2</u> |

Number TICs Found: 3

CONCENTRATION UNITS

| | CAS NO. | COMPOUND | RT | EST. CONC. | Q |
|----|------------|--------------------------|-------|------------|---|
| 1. | | Unknown | 4.363 | 3.17 | u |
| 2. | | Unknown | 4.623 | 9.62 | u |
| 3. | 49622-18-6 | Decane, 3,3,4-trimethyl- | 4.873 | 14.6 | u |

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1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

Lab Name: GCAL Sample ID: SKSWDEB1009
 Lab Code: LA024 Case No.: _____ Contract: _____
 SAS No.: _____ SDG No.: 204030804 Lab File ID: 2040319/S5001
 Matrix: Water Lab Sample ID: 20403080406
 Sample wt/vol: 1000 Units: mL Date Collected: 03/02/04 Time: 1240
 Level: (low/med) _____ Date Received: 03/06/04
 % Moisture: _____ decanted: (Y/N) _____ Date Extracted: 3/8/04
 GC Column: DB-5MS-30M ID: .25 (mm) Date Analyzed: 03/19/04 Time: 1513
 Concentrated Extract Volume: 1000 (µL) Dilution Factor: 1 Analyst: RLW
 Injection Volume: 1.0 (µL) Prep Method: _____
 GPC Cleanup: (Y/N) N pH: _____ Analytical Method: OLMO 4.2

Instrument ID: MSSV2

CONCENTRATION UNITS: ug/L

Prep Batch: 270522 Analytical Batch: 271099

| CAS NO. | COMPOUND | RESULT | Q | MDL | PQL |
|-----------|-----------------------------|--------|---|-------|------|
| 95-95-4 | 2,4,5-Trichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 88-06-2 | 2,4,6-Trichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 120-83-2 | 2,4-Dichlorophenol | 10.0 | U | 0.010 | 10.0 |
| 51-28-5 | 2,4-Dinitrophenol | 25.0 | U | 0.010 | 25.0 |
| 121-14-2 | 2,4-Dinitrotoluene | 10.0 | U | 0.010 | 10.0 |
| 806-20-2 | 2,6-Dinitrotoluene | 10.0 | U | 0.010 | 10.0 |
| 91-58-7 | 2-Chloronaphthalene | 10.0 | U | 0.010 | 10.0 |
| 95-57-8 | 2-Chlorophenol | 10.0 | U | 0.010 | 10.0 |
| 91-57-6 | 2-Methylnaphthalene | 10.0 | U | 0.010 | 10.0 |
| 88-74-4 | 2-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 88-75-5 | 2-Nitrophenol | 10.0 | U | 0.010 | 10.0 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 10.0 | U | 0.010 | 10.0 |
| 99-09-2 | 3-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 534-52-1 | 2-Methyl-4,6-dinitrophenol | 25.0 | U | 0.010 | 25.0 |
| 59-50-7 | 4-Chloro-3-methylphenol | 10.0 | U | 0.010 | 10.0 |
| 106-47-8 | 4-Chloroaniline | 10.0 | U | 0.010 | 10.0 |
| 7005-72-3 | 4-Chlorophenyl-phenylether | 10.0 | U | 0.010 | 10.0 |
| 106-44-5 | 4-Methylphenol (p-Cresol) | 10.0 | U | 0.010 | 10.0 |
| 83-32-9 | Acenaphthene | 10.0 | U | 0.010 | 10.0 |
| 208-96-8 | Acenaphthylene | 10.0 | U | 0.010 | 10.0 |
| 120-12-7 | Anthracene | 10.0 | U | 0.010 | 10.0 |
| 53-55-3 | Benzo(a)anthracene | 10.0 | U | 0.010 | 10.0 |
| 50-32-8 | Benzo(a)pyrene | 10.0 | U | 0.010 | 10.0 |
| 205-99-2 | Benzo(b)fluoranthene | 10.0 | U | 0.010 | 10.0 |
| 191-24-2 | Benzo(g,h,i)perylene | 10.0 | U | 0.010 | 10.0 |
| 207-08-9 | Benzo(k)fluoranthene | 10.0 | U | 0.010 | 10.0 |
| 111-91-1 | Bis(2-Chloroethoxy)methane | 10.0 | U | 0.010 | 10.0 |
| 111-44-4 | Bis(2-Chloroethyl)ether | 10.0 | U | 0.010 | 10.0 |
| 108-60-1 | bis(2-Chloroisopropyl)ether | 10.0 | U | 0.010 | 10.0 |
| 117-81-7 | bis(2-ethylhexyl)phthalate | 10.0 | U | 0.010 | 10.0 |

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SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

Lab Name: GCAL Sample ID: SKSWDEB1009
 Lab Code: LA024 Case No.: _____ Contract: _____
 SAS No.: _____ SOG No.: 204030804 Lab File ID: 2040319/S5001
 Matrix: Water Lab Sample ID: 20403080406
 Sample wt/vol: 1000 Units mL Date Collected: 03/02/04 Time: 1240
 Level: (low/med) _____ Date Received: 03/06/04
 % Moisture: _____ decanted: (Y/N) _____ Date Extracted: 3/8/04
 GC Column: DB-5MS-30M ID: 25 (mm) Date Analyzed: 03/19/04 Time: 1513
 Concentrated Extract Volume: 1000 (µL) Dilution Factor: 1 Analyst: RLW
 Injection Volume: 1.0 (µL) Prep Method: _____
 GPC Cleanup: (Y/N) N pH: _____ Analytical Method: OLMO 4.2

CONCENTRATION UNITS: ug/L

Instrument ID: MSSV2
 Prep Batch: 270522 Analytical Batch: 271099

| CAS NO. | COMPOUND | RESULT | Q | MDL | PQL |
|----------|---------------------------|--------|---|-------|------|
| 101-55-3 | 4-Bromophenyl-phenylether | 10.0 | U | 0.010 | 10.0 |
| 85-68-7 | Butylbenzylphthalate | 10.0 | U | 0.010 | 10.0 |
| 86-74-8 | Carbazole | 10.0 | U | 0.010 | 10.0 |
| 218-01-8 | Chrysene | 10.0 | U | 0.010 | 10.0 |
| 84-74-2 | Di-n-butylphthalate | 10.0 | U | 0.010 | 10.0 |
| 117-84-0 | Di-n-octylphthalate | 10.0 | U | 0.010 | 10.0 |
| 53-70-3 | Dibenz(a,h)anthracene | 10.0 | U | 0.010 | 10.0 |
| 132-64-9 | Dibenzofuran | 10.0 | U | 0.010 | 10.0 |
| 84-66-2 | Diethylphthalate | 10.0 | U | 0.010 | 10.0 |
| 131-11-3 | Dimethyl-phthalate | 10.0 | U | 0.010 | 10.0 |
| 105-67-9 | 2,4-Dimethylphenol | 10.0 | U | 0.010 | 10.0 |
| 206-44-0 | Fluoranthene | 10.0 | U | 0.010 | 10.0 |
| 86-73-7 | Fluorene | 10.0 | U | 0.010 | 10.0 |
| 118-74-1 | Hexachlorobenzene | 10.0 | U | 0.010 | 10.0 |
| 87-68-3 | Hexachlorobutadiene | 10.0 | U | 0.010 | 10.0 |
| 77-47-4 | Hexachlorocyclopentadiene | 10.0 | U | 0.010 | 10.0 |
| 67-72-1 | Hexachloroethane | 10.0 | U | 0.010 | 10.0 |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | 10.0 | U | 0.010 | 10.0 |
| 78-58-1 | Isophorone | 10.0 | U | 0.010 | 10.0 |
| 91-20-3 | Naphthalene | 10.0 | U | 0.010 | 10.0 |
| 100-01-6 | 4-Nitroaniline | 25.0 | U | 0.010 | 25.0 |
| 98-95-3 | Nitrobenzene | 10.0 | U | 0.010 | 10.0 |
| 100-02-7 | 4-Nitrophenol | 25.0 | U | 0.010 | 25.0 |
| 87-86-5 | Pentachlorophenol | 25.0 | U | 0.010 | 25.0 |
| 85-01-8 | Phenanthrene | 10.0 | U | 0.010 | 10.0 |
| 108-95-2 | Phenol | 10.0 | U | 0.010 | 10.0 |
| 129-00-0 | Pyrene | 10.0 | U | 0.010 | 10.0 |
| 621-64-7 | N-Nitroso-d-n-propylamine | 10.0 | U | 0.010 | 10.0 |
| 86-30-6 | N-Nitrosodiphenylamine | 10.0 | U | 0.010 | 10.0 |
| 95-48-7 | o-Cresol | 10.0 | U | 0.010 | 10.0 |

1F
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

Lab Name: GCAL Sample ID: SKSWDEB1009
Lab Code: LA024 2 Case No.: _____ Contract: _____
SAS No.: _____ SDG No.: 204030804 Lab File ID: 2040319/S5001
Matrix: Water Lab Sample ID: 20403080406
Sample wt/vol: _____ Units: _____ Date Collected: 03/02/04 Time: 1240
Level: (low/med) _____ Date Received: 03/06/04
% Moisture: not dec. _____ Date Extracted: 3/8/04
GC Column: DB-5MS-30M ID: .25 (mm) Date Analyzed: 03/19/04 Time: 1513
Concentrated Extract Volume: 1000 (µL) Dilution Factor: 1 Analyst: RLW
Injection Volume: 1.0 (µL) Prep Method: _____
GPC Cleanup: (Y/N) N pH: _____ Analytical Method: SW-846 8270C
Instrument ID: MSSV2

Number TICs Found : 3

CONCENTRATION UNITS:

| | CAS NO. | COMPOUND | RT | EST. CONC. | Q |
|----|------------|--------------------------|-------|------------|---|
| 1. | | Unknown | 4.366 | 3.57 | |
| 2. | | Unknown | 4.627 | 10.3 | |
| 3. | 49622-18-6 | Decane, 3,3,4-trimethyl- | 4.864 | 12 | |

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ORGANICS ANALYSIS DATA SHEET

Lab Name: GCAL Sample ID: SKSWD031009
 Lab Code: LA024 Case No.: _____ Contract: _____
 Matrix: Water SAS No.: _____ SDG No.: 204030804
 Sample wt/vol: 1000 Units: mL Lab Sample ID: 20403080401
 Level: (low/med) _____ Date Collected: 03/02/04 Time: 1055
 % Moisture: _____ deaerated (Y/N) _____ Date Received: 03/06/04
 GC Column: DB-608-30M ID: .53 (mm) Date Extracted: 3/8/04 *sho for*
 Concentrated Extract Volume: 1000 (µL) Date Analyzed: 03/31/04 Time: 0540
 Soil Aliquot Volume: _____ (µL) Dilution Factor: 1 Analyst: TLS
 Injection Volume: 1 (µL) Prep Method: _____
 GPC Cleanup: (Y/N) N pH: _____ Analytical Method: OLMO 4.2
 Prep Batch: 270511 Analytical Batch: 271554 Sulfur Cleanup: (Y/N) N Instrument ID: GCS6A
 Lab File ID: 2040330/SV6023

CONCENTRATION UNITS: µg/L

CAS NO. COMPOUND RESULT Q MDL RL

| | | | | | |
|------------|---------------------|-------|---|---------|-------|
| 72-54-8 | 4,4'-DDD | 0.100 | U | 0.00010 | 0.100 |
| 72-55-9 | 4,4'-DDE | 0.100 | U | 0.00010 | 0.100 |
| 50-29-3 | 4,4'-DDT | 0.100 | U | 0.00010 | 0.100 |
| 309-00-2 | Aldrin | 0.050 | U | 0.00010 | 0.050 |
| 12674-11-2 | Aroclor-1016 | 1.00 | U | 0.00010 | 1.00 |
| 11104-28-2 | Aroclor-1221 | 2.00 | U | 0.00010 | 2.00 |
| 11141-16-5 | Aroclor-1232 | 1.00 | U | 0.00010 | 1.00 |
| 53469-21-9 | Aroclor-1242 | 1.00 | U | 0.00010 | 1.00 |
| 12672-29-6 | Aroclor-1248 | 1.00 | U | 0.00010 | 1.00 |
| 11097-69-1 | Aroclor-1254 | 1.00 | U | 0.00010 | 1.00 |
| 11096-82-5 | Aroclor-1260 | 1.00 | U | 0.00010 | 1.00 |
| 60-57-1 | Dieldrin | 0.100 | U | 0.00010 | 0.100 |
| 959-98-8 | Endosulfan I | 0.050 | U | 0.00010 | 0.050 |
| 33213-65-9 | Endosulfan II | 0.100 | U | 0.00010 | 0.100 |
| 1031-07-8 | Endosulfan sulfate | 0.100 | U | 0.00010 | 0.100 |
| 72-20-8 | Endrin | 0.100 | U | 0.00010 | 0.100 |
| 7421-93-4 | Endrin aldehyde | 0.100 | U | 0.00010 | 0.100 |
| 53494-70-5 | Endrin ketone | 0.100 | U | 0.00010 | 0.100 |
| 76-44-8 | Heptachlor | 0.050 | U | 0.00010 | 0.050 |
| 1024-57-3 | Heptachlor epoxide | 0.050 | U | 0.00010 | 0.050 |
| 72-43-5 | Methoxychlor | 0.500 | U | 0.00010 | 0.500 |
| 8001-35-2 | Toxaphene | 5.00 | U | 0.00010 | 5.00 |
| 319-84-6 | alpha-BHC | 0.050 | U | 0.00010 | 0.050 |
| 5103-71-9 | alpha-Chlordane | 0.050 | U | 0.00010 | 0.050 |
| 319-85-7 | beta-BHC | 0.050 | U | 0.00010 | 0.050 |
| 319-86-8 | delta-BHC | 0.050 | U | 0.00010 | 0.050 |
| 58-89-9 | gamma-BHC (Lindane) | 0.050 | U | 0.00010 | 0.050 |
| 5103-74-2 | gamma-Chlordane | 0.050 | U | 0.00010 | 0.050 |

1D
ORGANICS ANALYSIS DATA SHEET

Lab Name: GCAL Sample ID: SKSWD03D1009
 Lab Code: LA024 Case No.: _____ Contract: _____
 Matrix: Water SAS No.: _____ SDG No.: 204030804
 Sample wt/vol: 1000 Units: mL Lab Sample ID: 20403080402
 Level: (low/med) _____ Date Collected: 03/02/04 Time: 1118
 % Moisture: _____ decanted: (Y/N) _____ Date Received: 03/06/04
 GC Column: DB-608-30M ID: .53 (mm) Date Extracted: 3/8/04 *shahar msa*
 Concentrated Extract Volume: 1000 (µL) Date Analyzed: 03/31/04 Time: 0608
 Soil Aliquot Volume: _____ (µL) Dilution Factor: 1 Analyst: TLS
 Injection Volume: 1 (µL) Prep Method: _____
 GPC Cleanup: (Y/N) N pH: _____ Analytical Method: OLMO 4.2
 Prep Batch: 270511 Analytical Batch: 271554 Sulfur Cleanup: (Y/N) N Instrument ID: GCS6A
 Lab File ID: 2040330/SV6024

CONCENTRATION UNITS: ug/L

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|------------|---------------------|--------|---|---------|-------|
| 72-54-8 | 4,4'-DDD | 0.100 | U | 0.00010 | 0.100 |
| 72-55-9 | 4,4'-DDE | 0.100 | U | 0.00010 | 0.100 |
| 50-29-3 | 4,4'-DDT | 0.100 | U | 0.00010 | 0.100 |
| 309-00-2 | Aldrin | 0.050 | U | 0.00010 | 0.050 |
| 12674-11-2 | Aroclor-1016 | 1.00 | U | 0.00010 | 1.00 |
| 11104-28-2 | Aroclor-1221 | 2.00 | U | 0.00010 | 2.00 |
| 11141-16-5 | Aroclor-1232 | 1.00 | U | 0.00010 | 1.00 |
| 53469-21-9 | Aroclor-1242 | 1.00 | U | 0.00010 | 1.00 |
| 12672-29-6 | Aroclor-1248 | 1.00 | U | 0.00010 | 1.00 |
| 11097-69-1 | Aroclor-1254 | 1.00 | U | 0.00010 | 1.00 |
| 11096-82-5 | Aroclor-1260 | 1.00 | U | 0.00010 | 1.00 |
| 60-57-1 | Dieldrin | 0.100 | U | 0.00010 | 0.100 |
| 959-98-8 | Endosulfan I | 0.050 | U | 0.00010 | 0.050 |
| 33213-65-9 | Endosulfan II | 0.100 | U | 0.00010 | 0.100 |
| 1031-07-8 | Endosulfan sulfate | 0.100 | U | 0.00010 | 0.100 |
| 72-20-8 | Endrin | 0.100 | U | 0.00010 | 0.100 |
| 7421-93-4 | Endrin aldehyde | 0.100 | U | 0.00010 | 0.100 |
| 53494-70-5 | Endrin ketone | 0.100 | U | 0.00010 | 0.100 |
| 76-44-8 | Heptachlor | 0.050 | U | 0.00010 | 0.050 |
| 1024-57-3 | Heptachlor epoxide | 0.050 | U | 0.00010 | 0.050 |
| 72-43-5 | Methoxychlor | 0.500 | U | 0.00010 | 0.500 |
| 8001-35-2 | Toxaphene | 5.00 | U | 0.00010 | 5.00 |
| 319-84-6 | alpha-BHC | 0.050 | U | 0.00010 | 0.050 |
| 5103-71-9 | alpha-Chlordane | 0.050 | U | 0.00010 | 0.050 |
| 319-85-7 | beta-BHC | 0.050 | U | 0.00010 | 0.050 |
| 319-86-8 | delta-BHC | 0.050 | U | 0.00010 | 0.050 |
| 58-89-9 | gamma-BHC (Lindane) | 0.050 | U | 0.00010 | 0.050 |
| 5103-74-2 | gamma-Chlordane | 0.050 | U | 0.00010 | 0.050 |

ORGANICS ANALYSIS DATA SHEET

Lab Name: GCAL Sample ID: SKS.VDEB1009
 Lab Code: LA024 Case No.: _____ Contract: _____
 Matrix: Water SAS No.: _____ SDG No.: 204030804
 Sample wt/vol: 1000 Units mL Lab Sample ID: 20403080406
 Level: (low/med) _____ Date Collected: 03/02/04 Time: 1240
 % Moisture: _____ decanted: (Y/N) _____ Date Received: 03/06/04
 GC Column: DB-608-30M ID: .53 (mm) Date Extracted: 3/2/04
 Concentrated Extract Volume: 1000 (µL) Date Analyzed: 03/31/04 Time: 0733
 Soil Aliquot Volume: _____ (µL) Dilution Factor: 1 Analyst: TLS
 Injection Volume: 1 (µL) Prep Method: _____
 GPC Cleanup: (Y/N) N pH: _____ Analytical Method: OLMO 4.2
 Prep Batch: 270511 Analytical Batch: 271554 Surfer Cleanup (Y/N) N Instrument ID: GCS6A

CONCENTRATION UNITS: µg/L

Lab File ID: 2040330/SV6027

| CAS NO. | COMPOUND | RESULT | Q | MDL | RL |
|------------|---------------------|--------|---|---------|-------|
| 72-54-8 | 4,4'-DDD | 0.100 | U | 0.00010 | 0.100 |
| 72-55-9 | 4,4'-DDE | 0.100 | U | 0.00010 | 0.100 |
| 50-29-3 | 4,4'-DDT | 0.100 | U | 0.00010 | 0.100 |
| 309-00-2 | Aldrin | 0.050 | U | 0.00010 | 0.050 |
| 12674-11-2 | Aroclor-1016 | 1.00 | U | 0.00010 | 1.00 |
| 11104-28-2 | Aroclor-1221 | 2.00 | U | 0.00010 | 2.00 |
| 11141-16-5 | Aroclor-1232 | 1.00 | U | 0.00010 | 1.00 |
| 53469-21-9 | Aroclor-1242 | 1.00 | U | 0.00010 | 1.00 |
| 12672-29-6 | Aroclor-1248 | 1.00 | U | 0.00010 | 1.00 |
| 11097-69-1 | Aroclor-1254 | 1.00 | U | 0.00010 | 1.00 |
| 11096-82-5 | Aroclor-1260 | 1.00 | U | 0.00010 | 1.00 |
| 60-57-1 | Dieldrin | 0.100 | U | 0.00010 | 0.100 |
| 959-98-8 | Endosulfan I | 0.050 | U | 0.00010 | 0.050 |
| 33213-65-9 | Endosulfan II | 0.100 | U | 0.00010 | 0.100 |
| 1031-07-8 | Endosulfan sulfate | 0.100 | U | 0.00010 | 0.100 |
| 72-20-8 | Endrin | 0.100 | U | 0.00010 | 0.100 |
| 7421-93-4 | Endrin aldehyde | 0.100 | U | 0.00010 | 0.100 |
| 53494-70-5 | Endrin ketone | 0.100 | U | 0.00010 | 0.100 |
| 76-44-8 | Heptachlor | 0.050 | U | 0.00010 | 0.050 |
| 1024-57-3 | Heptachlor epoxide | 0.050 | U | 0.00010 | 0.050 |
| 72-43-5 | Methoxychlor | 0.500 | U | 0.00010 | 0.500 |
| 8001-35-2 | Toxaphene | 5.00 | U | 0.00010 | 5.00 |
| 319-84-6 | alpha-BHC | 0.050 | U | 0.00010 | 0.050 |
| 5103-71-9 | alpha-Chlordane | 0.050 | U | 0.00010 | 0.050 |
| 319-85-7 | beta-BHC | 0.050 | U | 0.00010 | 0.050 |
| 319-86-8 | delta-BHC | 0.050 | U | 0.00010 | 0.050 |
| 58-89-9 | gamma-BHC (Lincane) | 0.050 | U | 0.00010 | 0.050 |
| 5103-74-2 | gamma-Chlordane | 0.050 | U | 0.00010 | 0.050 |

U.S. EPA - CLP
1
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKSWD031009

Lab Name: GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204030804
Matrix: (soil / water) Water Lab Sample ID: 20403080401
Level: (low / med) _____ Date Received: 03/06/04
% Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7440-36-0 | Antimony | 5.2 | B | | P |
| 7440-38-2 | Arsenic | 2.9 | U | | P |
| 7440-39-3 | Barium | 40.0 | B | | P |
| 7440-41-7 | Beryllium | 0.2 | B | | P |
| 7440-43-9 | Cadmium | 0.2 | U | | P |
| 7440-47-3 | Chromium | 1.4 | B | | P |
| 7440-50-8 | Copper | 11.0 | B | | P |
| 7439-89-6 | Iron | 2200 | | | P |
| 7439-92-1 | Lead | 1.5 | U | | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 2.5 | B | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-66-6 | Zinc | 14.6 | B | | P |
| 57-12-5 | Cyanide | 0.8 | B | | AS |

Si nloy
msk

Color Before: COLORLESS Clarity Before: CLEAR Texture: _____
Color After: COLORLESS Clarity After: CLEAR Artifacts: _____
Comments:

U.S. EPA - OLP
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKSWD03D1009

Lab Name: GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204030804
Matrix (soil / water) Water Lab Sample ID 20403080402
Level: (low / med) _____ Date Received: 03/06/04
% Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7440-36-0 | Antimony | 3.7 | L | | P |
| 7440-38-2 | Arsenic | 5.3 | B | | P |
| 7440-35-3 | Barium | 39.1 | B | | P |
| 7440-41-7 | Beryllium | 0.2 | B | | P |
| 7440-43-9 | Cadmium | 0.2 | L | | P |
| 7440-47-3 | Chromium | 1.3 | B | | P |
| 7440-50-8 | Copper | 4.5 | B | | P |
| 7439-89-6 | Iron | 2220 | | | P |
| 7439-92-1 | Lead | 1.5 | L | | P |
| 7439-97-6 | Mercury | 0.1 | L | | AV |
| 7440-02-0 | Nickel | 2.8 | B | | P |
| 7782-49-2 | Selenium | 4.4 | L | N | P |
| 7440-22-4 | Silver | 0.4 | L | | P |
| 7440-28-0 | Thallium | 2.6 | L | | P |
| 7440-66-6 | Zinc | 7.2 | B | | P |
| 57-12-5 | Cyanide | 2.5 | B | | AS |

slightly
pink

Color Before: COLORLESS Clarity Before: CLEAR Texture: _____
Color After: COLORLESS Clarity After: CLEAR Artifacts: _____
Comments: _____

U.S. EPA - CLP
1
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKSWDEB1009

Lab Name: GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204030804
Matrix: (soil / water) Water Lab Sample ID: 20403080406
Level: (low / med) _____ Date Received: 03/06/04
% Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7440-36-0 | Antimony | 3.7 | U | | P |
| 7440-38-2 | Arsenic | 2.9 | U | | P |
| 7440-39-3 | Barium | 0.3 | U | | P |
| 7440-41-7 | Beryllium | 0.1 | U | | P |
| 7440-43-9 | Cadmium | 0.2 | U | | P |
| 7440-47-3 | Chromium | 0.8 | U | | P |
| 7440-50-8 | Copper | 3.5 | B | | P |
| 7439-89-6 | Iron | 14.1 | U | | P |
| 7439-92-1 | Lead | 1.5 | U | | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 0.7 | U | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-66-6 | Zinc | 3.7 | B | | P |
| 57-12-5 | Cyanide | 1.9 | B | | AS |

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5/11/04
msc

Color Before: COLORLESS

Clarity Before: CLEAR

Texture: _____

Color After: COLORLESS

Clarity After: CLEAR

Artifacts: _____

Comments:

U.S. EPA - DLP
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKSWD031009 (DISS)

Lab Name: GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204030804
Matrix (soil / water): Water Lab Sample ID: 20403080408
Level: (low / med) _____ Date Received: 03/06/04
% Solids: _____

Concentration Units (ug/L or mg/kg dry weight): ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7440-36-0 | Antimony | 3.7 | L | | P |
| 7440-38-2 | Arsenic | 5.3 | B | | P |
| 7440-39-3 | Barium | 29.8 | B | | P |
| 7440-41-7 | Beryllium | 0.2 | B | | P |
| 7440-43-9 | Cadmium | 0.2 | U | | P |
| 7440-47-3 | Chromium | 0.8 | L | | P |
| 7440-50-8 | Copper | 4.6 | B | | P |
| 7439-89-6 | Iron | 17.2 | B | | P |
| 7439-92-1 | Lead | 1.5 | U | | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 1.4 | B | | P |
| 7782-49-2 | Selenium | 4.4 | L | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-28-0 | Thallium | 2.6 | L | | P |
| 7440-66-6 | Zinc | 0.6 | L | | P |

J

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5/11/04
msl

Color Before: COLORLESS Clarity Before: CLEAR Texture: _____
Color After: COLORLESS Clarity After: CLEAR Artifacts: _____
Comments: _____

U.S. EPA - CLP
1
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKSWD03D1009 (DISS)

Lab Name: GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204030804
Matrix: (soil / water) Water Lab Sample ID: 20403080409
Level: (low / med) _____ Date Received: 03/06/04
% Solids: _____
Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7440-36-0 | Antimony | 3.7 | U | | P |
| 7440-38-2 | Arsenic | 4.1 | B | | P |
| 7440-39-3 | Barium | 29.8 | B | | P |
| 7440-41-7 | Beryllium | 0.1 | B | | P |
| 7440-43-9 | Cadmium | 0.2 | U | | P |
| 7440-47-3 | Chromium | 0.8 | U | | P |
| 7440-50-8 | Copper | 1.2 | U | | P |
| 7439-89-6 | Iron | 27.7 | B | | P |
| 7439-92-1 | Lead | 1.5 | U | | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 1.3 | B | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-66-6 | Zinc | 0.6 | U | | P |

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5/11/04
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Color Before: COLORLESS Clarity Before: CLEAR Texture: _____
Color After: COLORLESS Clarity After: CLEAR Artifacts: _____
Comments: _____

U.S. EPA - OLP
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKSWDEB1009(DISS)

Lab Name: GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204030804
Matrix (soil / water): Water Lab Sample ID: 20403080412
Level (low / med): _____ Date Received: 03/06/04
% Solids: _____

Concentration Units (ug/L or mg/kg dry weight): ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7440-36-0 | Antimony | 3.7 | U | | P |
| 7440-38-2 | Arsenic | 2.9 | U | | P |
| 7440-39-3 | Barium | 0.3 | U | | P |
| 7440-41-7 | Beryllium | 0.1 | U | | P |
| 7440-43-8 | Cadmium | 0.2 | U | | P |
| 7440-47-3 | Chromium | 0.8 | U | | P |
| 7440-50-8 | Copper | 1.2 | U | | P |
| 7439-89-8 | Iron | 22.6 | B | | P |
| 7439-92-1 | Lead | 1.5 | U | | P |
| 7439-97-8 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 1.5 | B | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-66-6 | Zinc | 3.9 | B | | P |

* 43

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Color Before: COLORLESS Clarity Before: CLEAR Texture: _____
Color After: COLORLESS Clarity After: CLEAR Artifacts: _____
Comments: _____

CHAIN OF CUSTODY RECORD

Lab use only

Earth Tech

Client Name

4342

Client #

204030804

Workorder #

3/19/04

Due Date

Report to:

Client: *Earth Tech*

Address: *200 Vine Street
Wilders, KY 41076*

Contact: *Pat Higgins*

Phone: *859-442-2300*

Fax: *859-442-2311*

Bill to:

Client: _____

Address: _____

Contact: *same*

Phone: _____

Fax: _____

Analytical Requests & Method

*Volatiles
Semi Volatiles
Pesticides
PCBs
total metals
dissolved metals
cyanide*

Lab use only:

Custody Seal

used ☒ yes ☐ no

in tact ☒ yes ☐ no

Temperature °C *4*

P.O. Number

54280

Project Name/Number

Skinner Landfill - 1 Qtr. 2004

Sampled By:

Pat Higgins / Chris Cox

| Matrix ¹ | Date | Time (2400) | Comp | Gr | Sample Description | Preservatives | No Containers | Volatiles | Semi Volatiles | Pesticides | PCBs | total metals | dissolved metals | Cyanide | Remarks | Lab ID |
|---------------------|--------|-------------|------|----|--------------------|---------------|---------------|-----------|----------------|------------|------|--------------|------------------|---------|----------------------|--------|
| W | 3/2/04 | 1055 | | X | SKSWD031009 | HCL | 3 | X | X | X | X | X | X | X | 8 Refer to Table 7 | -1 |
| | | 1118 | | | SKSWD03D1009 | HCL | 3 | X | X | X | X | X | X | X | 9 (TCL) and Table 8 | 2 |
| | | 1145 | | | SKSWD03MS 1009 | HCL | 7 | X | | | | | | | 10 (TAL) of final | 3 |
| | | 1208 | | | SKSWD03 MSP 1009 | HCL | 7 | X | | | | | | | 11 O&M plan for dup. | 5 |
| | | 1240 | | | SKSWD03B 1009 | Various | 10 | X | X | X | X | X | X | X | 12 list of analytes. | 6 |
| | | | | | Trip Blank | HCL | 3 | X | | | | | | | | 7 |

Turn Around Time: ☐ 24-48 hrs. ☐ 3 days ☐ 1 week ☒ Standard ☐ Other _____

Relinquished by: (Signature)

Pat Higgins

Received by: (Signature)

FedX

Date:

Time:

Note:

Relinquished by: (Signature)

FedEX 8425 4530 7654

Received by: (Signature)

Don Miller

Date:

Time:

3-6-04

1030

Relinquished by: (Signature)

Don Miller

Received by: (Signature)

Cooper

Date:

Time:

3-6-04

1150

By submitting these samples, you agree to the terms and conditions contained in our most recent schedule of services.

CHAIN OF CUSTODY RECORD

Lab use only

Earth Tech

Client Name

4342

Client #

704030804

Workorder #

5/19/04

Run Date

Report to:

Client: Earth Tech
Address: 200 Vine Street
Wilders, KY 41076
Contact: Pat Higgins
Phone: 859 442-2700
Fax: 859 442-2111

Bill to:

Client:
Address:
Contact: SANG
Phone:
Fax:

P.O. Number

54280

Project Name/Number

Shiner Landfill- 1 Qtr. 2004

Sampled By:

Pat Higgins/Chris Cox

| Matrix | Date | Time (2400) | Sample Description | Preservatives | No. Containers |
|--------|--------|-------------|--------------------|---------------|----------------|
| W | 3/2/04 | 1055 | X Sk SWD03 1009 | Varia | 7 |
| W | 3/2/04 | 1118 | X Sk SWD03D 1009 | Varia | 7 |

Analytical Requests & Method

| Volatiles | Semi Volatiles | PCRS | Pesticides | Total Metals | Dissolved Metals | Cyanide |
|-----------|----------------|------|------------|--------------|------------------|---------|
| X | X | X | X | X | X | X |
| X | X | X | X | X | X | X |

Lab use only:

Custody Seal

used ☒ yes ☐ no

In tact ☒ yes ☐ no

Temperature °C

Remarks:

8 Refer to
9 Table 7
(TCL) and
Table 8
(TAL) of
Final O&M
Plan for
list of
analytes

Lab ID

3/1
-1
2

Turn Around Time: ☐ 24-48 hrs. ☐ 3 days ☐ 1 week ☒ Standard ☐ Other

Relinquished by: (Signature)

Pat Higgins

Relinquished by: (Signature)

Fedex 9425 9530 7656

Relinquished by: (Signature)

Don Miller

Received by: (Signature)

FED X

Received by: (Signature)

Don Miller

Received by: (Signature)

Cooper

Date:

Time:

Date:

Time:

Date:

Time:

Note:

By submitting these samples, you agree to the terms and conditions contained in our most recent schedule of services

CHAIN OF CUSTODY RECORD

| | | | | |
|--------------|-------------|----------|-------------|----------|
| Lab use only | Earth Tech | 4342 | 204030804 | 3/19/04 |
| | Client Name | Client # | Workorder # | Due Date |

| Report to: | | Bill to: | | Analytical Requests & Method | | | | | | | | | | Lab use only: | | | | |
|--|--------|--|-----|---|----------------------|---------------|---------------|---|---|---|---|---|---|---|-----|--|--|--------|
| Client: Earth Tech | | Client: | | Volatiles Semi Volatiles Pesticides PCB's Total Metals Dissolved Metals Cyanide | | | | | | | | | | Custody Seal | | | | |
| Address: 200 Vine Street Wilder, LA 70766 | | Address: | | | | | | | | | | | | used <input checked="" type="checkbox"/> yes <input type="checkbox"/> no | | | | |
| Contact: Pat Higgins | | Contact: Same | | | | | | | | | | | | in tact <input checked="" type="checkbox"/> yes <input type="checkbox"/> no | | | | |
| Phone: 854 442 2700 | | Phone: | | | | | | | | | | | | Temperature °C 4 | | | | |
| Fax: 854 442 2311 | | Fax: | | | | | | | | | | | | | | | | |
| P.O. Number 54280 | | Project Name/Number Shinner Landfill - 1 Qtr. 2004 | | | | | | | | | | | | | | | | |
| Sampled By: Pat Higgins / Chris Cox | | | | | | | | | | | | | | | | | | |
| Matrix | Date | Time (2400) | COB | G | Sample Description | Preservatives | No Containers | | | | | | | | | | | Lab ID |
| W | 3/2/04 | 1145 | | X | Sh SW D 03 MS 1004 | Various | 7 | X | X | X | X | X | X | 10 | 3/8 | | | |
| W | 3/2/04 | 1208 | | X | Sh SW D 03 MS D 1004 | Various | 7 | X | X | X | X | X | X | 11 MOD-4 | 3 | | | |
| Remarks: Refer to Table 7 and Table 8 of Final O&M Plan for list of analytes | | | | | | | | | | | | | | | | | | |

Turn Around Time: ☐ 24-48 hrs. ☐ 3 days ☐ 1 week ☒ Standard ☐ Other

| | | | | |
|--|--|-------|-------|-------|
| Relinquished by: (Signature) Pat Higgins | Received by: (Signature) FED X | Date: | Time: | Note: |
| Relinquished by: (Signature) FED X 0415 2530 7645 | Received by: (Signature) Don Palmer | Date: | Time: | |
| Relinquished by: (Signature) Don Palmer | Received by: (Signature) Cooler | Date: | Time: | |

By submitting these samples, you agree to the terms and conditions contained in our most recent schedule of services.

DATA VALIDATION REPORT
FOR
SKINNER LANDFILL SITE
EARTH TECH: PROJECT NUMBER 54280
LABORATORY REPORT NUMBER 204031909
PROJECT MANAGER: Ron Rolker
Date: May 11, 2004
Revised Report Dated: September 29, 2005
Data Validator: Mark Kromis

APPENDIX C LIST OF ACRONYMS

| | |
|--------|--|
| BFB | Bromofluorobenzene |
| CC | Continuing Calibration |
| CCV | Continuing Calibration Verification |
| CCB | Continuing Calibration Blanks |
| CLP | Contract Laboratory Program |
| CRDL | Contract Required Detection Limit |
| DFTPP | Decafluorotriphenylphosphine |
| GC/MS | Gas Chromatograph/Mass Spectrometer |
| IC | Initial Calibration |
| ICB | Initial Calibration Blank |
| IDL | Instrument Detection Limit |
| ICP | Inductively Coupled Plasma |
| ICS | Interference Check Sample |
| ICV | Initial Calibration Verification |
| ILM | Inorganic Analysis Multi-Media Multi-Concentration |
| INDAM | Individual A Mixture |
| INDBM | Individual B Mixture |
| mg/L | milligrams per liter |
| MS/MSD | Matrix Spike/Matrix Spike Duplicate |
| OLC | Organic Analysis Low Concentration |
| OLM | Organic Analysis Multi-Media Multi-Concentration |
| %D | Percent Difference |
| % RSD | Percent Relative Standard Deviation |
| PB | Preparation Blanks |
| QC | Quality Control |
| RF | Response Factor |
| RPD | Relative Percent Difference |
| RRF | Relative Response Factor |
| SDG | Sample Delivery Group |
| SOW | Statement of Work |
| µg/L | micrograms per liter |
| US EPA | United States Environmental Protection Agency |
| VOC | Volatile Organic Compounds |
| VTSR | Validated Time of Sample Receipt |

**DATA VALIDATION SUMMARY – SAMPLE DELIVERY GROUP 204031909
INORGANICS**

Validation of the inorganics data, as prepared by Gulf Coast Analytical Laboratories (GCAL) for the samples collected from the Skinner Landfill site in March 2004, was conducted by Earth Tech using the National Functional Guidelines for Inorganic Data Review, (US EPA, February, 1994), as appropriate. The results were reported by GCAL under Sample Delivery Group (SDG) 204031909.

| GCAL # | Sample Description |
|---------------|---------------------------|
| 20403190901 | SKGW581009 |
| 20403190902 | SKGW58DUP1009 |
| 20403190903 | SKGW631009 |
| 20403190904 | SKGW611009 |
| 20403190905 | SKGW61MS1009 |
| 20403190907 | SKGW61DUP1009 |
| 20403190908 | SKGWFB1009 |
| 20403190910 | SKGW581009 (DISS) |
| 20403190911 | SKGW58DUP1009 (DISS) |
| 20403190912 | SKGW631009 (DISS) |
| 20403190913 | SKGW611009 (DISS) |
| 20403190914 | SKGW61MS1009 (DISS) |
| 20403190915 | SKGW61DUP1009 (DISS) |
| 20403190916 | SKGWFB1009 (DISS) |
| 20403220801 | SKGW06R1009 |
| 20403220802 | SKGW07R1009 |
| 20403220803 | SKGW591009 |
| 20403220804 | SKGW601009 |
| 20403220805 | SKGW62A1009 |
| 20403220806 | SKGW641009 |
| 20403220807 | SKGW651009 |
| 20403220808 | SKGW06R1009 (DISS) |
| 20403220809 | SKGW07R1009 (DISS) |
| 20403220810 | SKGW591009 (DISS) |
| 20403220811 | SKGW601009 (DISS) |
| 20403220812 | SKGW62A1009 (DISS) |
| 20403220813 | SKGW641009 (DISS) |

INTRODUCTION

Analyses of metals were performed according to Contract Laboratory Program (CLP)-Inorganic Analysis Multi-media Multi-concentration ILM04.1 Statement of Work (SOW). Results of the sample analyses are reported by the laboratory as either qualified or unqualified. Unqualified results mean that the reported values maybe used without reservation. Various qualifier codes are used by the laboratory to denote specific information regarding the analytical results.

The data validation process is intended to evaluate the data on a technical basis. The data package also was subjected to an internal laboratory quality review prior to submission to Earth Tech for data validation.

During the validation process, laboratory-qualified and unqualified data are verified against all available supporting documentation. Based on this evaluation, qualifier codes may be added, deleted or modified by the data user. Final results are therefore, either qualified or unqualified. Validator-qualified results are annotated with the following codes in accordance with the Functional Guidelines:

- U The constituent was analyzed for, but was not detected above the level of the associated analytical reporting limit. The associated value is either the sample quantitation limit or the sample detection limit.
- J The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample.
- UJ The analyte was not detected above the reported sample quantitation limit. However, the reported quantitation limit is approximate and may or may not represent the actual limit of quantitation necessary to accurately and precisely measure the analyte in the sample.
- R The sample results are rejected due to serious deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte cannot be verified.

Details of the inorganics data validation findings and conclusions are provided in the following sections of this report:

1. Holding Times
2. Calibration
 - A. Initial Calibration (IC)
 - B. Continuing Calibration (CC)
3. Blanks
4. Inductively Coupled Plasma (ICP) Interference Check Sample
5. Laboratory Control Sample (LCS)
6. Duplicate Analysis

7. Spike Sample Analysis
8. ICP Serial Dilution
9. System Performance
10. Documentation
11. Overall Assessment

1. HOLDING TIMES

All samples for inorganics analyses were analyzed within the 180-day holding time for preserved aqueous samples. Mercury analyses were conducted within the 28-day holding time for aqueous samples undergoing CLP protocol. Cyanide analyses were conducted within the 14-day holding time. The cooler temperature upon receipt at the laboratory was within the recommended temperature of 4°C +/- 2°C.

2. CALIBRATION

A. Initial Calibration

The percent recoveries for the Initial Calibration Verification (ICV) standard were within Quality Control (QC) limits for all constituents.

B. Continuing Calibration

The percent recoveries for the Continuing Calibration Verification (CCV) standard were within QC limits for all constituents.

3. BLANKS

The Initial Calibration Blank (ICB), Continuing Calibration Blanks (CCB) and Preparation Blanks (PB) were analyzed at the appropriate frequencies. No analytes were detected in the ICB, CCB, PB and Field blanks above the corresponding Contract Required Detection Limit (CRDL).

4. ICP INTERFERENCE CHECK SAMPLE

Results for the ICP analysis of the Interference Check Sample (ICS) solution AB were within 20% of the true value.

5. LABORATORY CONTROL SAMPLES

Recoveries were within the contract (90-120%) for all constituents.

6. DUPLICATE ANALYSIS

The Relative Percent Difference (RPD) between the sample and duplicate results were within the acceptance criteria for all target compounds.

7. SPIKE SAMPLE ANALYSIS

The laboratory used sample SKGW611009 for the matrix spike sample. The MS percent recoveries were within the acceptance criteria (75%-125%) with the exception of Selenium in the total (45%) and dissolved (73%) fractions. As per the National Functional Guidelines: if the percent recovery is greater than 30% and less than 74% qualify detected results for that analyte with "J" and non-detected results with "UJ".

8. ICP SERIAL DILUTION

As noted in the National Functional Guidelines: If the analyte concentration is at least 50 times above the IDL, its serial dilution analysis must then agree within 10% of the original determination after corrected for dilution. The serial dilution is performed to determine whether any significant chemical or physical interference's exist due to matrix effects. The serial dilution percent differences were within the acceptance criteria for all target analytes with the exception of Manganese associated with the dissolved fraction.

9. SYSTEM PERFORMANCE

The analytical system appears to have been working well at the time of these analyses, based on the evaluation of the raw data.

10. DOCUMENTATION

GCAL qualified the total metal results for Aluminum, Copper, Lead, and Zinc with an "E" qualifier on the Form 1's to indicate that the %Difference (%D) between the original results and its serial dilution result exceeded the control limit. GCAL qualified the dissolved metal results for Iron with an "E" qualifier on the Form 1's to indicate that the %Difference (%D) between the original results and its serial dilution result exceeded the control limit. The %Differences were actually within the control limit therefore the data validator crossed out the (E) with a single line and dated and initialed the bottom of the report.

11. OVERALL ASSESSMENT

The percent recoveries for Arsenic in the Contract Required Detection Limit (CRDL) standards were 123%, 106%, and 120%.

The percent recoveries for Copper in the Contract Required Detection Limit (CRDL) standards were 84, 81%, and 79%.

The percent recoveries for Selenium in the Contract Required Detection Limit (CRDL) standards were 122%, 100%, and 139%. The Selenium results were previously qualified under Section 7-titled "Spike Sample Analysis".

The percent recoveries for Thallium in the Contract Required Detection Limit (CRDL) standards were 125.8%, 146.7%, and 130.4%.

The percent recoveries for Zinc in the Contract Required Detection Limit (CRDL) standards were 73.9%, 71.5%, and 68.9%.

If the CRDL is greater than 120% then detected results greater than the IDL but less than two times the CRDL are qualified as estimated with "J". If the CRDL is below 80% then detected results are qualified as estimated with "J" and the non-detected results were qualified with "UJ".

The results are acceptable with the validator-added qualifiers.

**DATA VALIDATION SUMMARY – SAMPLE DELIVERY GROUP 204031909
SEMIVOLATILE ORGANICS**

Validation of the Gas Chromatograph/Mass Spectrometer (GC/MS) semi-volatile organics data, as prepared by Gulf Coast Analytical Laboratories (GCAL) for the samples collected from the Skinner Landfill site in March 2004, was conducted by Earth Tech using the National Functional Guidelines for Organic Data Review, (US EPA, October, 1999) as appropriate. The results were reported by GCAL under SDG 204031909.

| GCAL # | Sample Description |
|---------------|---------------------------|
| 20403190901 | SKGW581009 |
| 20403190902 | SKGW58DUP1009 |
| 20403190903 | SKGW631009 |
| 20403190904 | SKGW611009 |
| 20403190905 | SKGW61MS1009 |
| 20403190906 | SKGW61MSD1009 |
| 20403190908 | SKGWFB1009 |
| 20403220801 | SKGW06R1009 |
| 20403220802 | SKGW07R1009 |
| 20403220803 | SKGW591009 |
| 20403220804 | SKGW601009 |
| 20403220805 | SKGW62A1009 |
| 20403220806 | SKGW641009 |
| 20403220807 | SKGW651009 |

INTRODUCTION

Analyses were performed according to CLP-Organic Analysis Multi-Media, Multi-Concentration OLM04.2 SOW. Results of the sample analyses are reported by the laboratory as either qualified or unqualified. Unqualified results mean that the reported values may be used without reservation. Various data qualifier codes are used by the laboratory to denote specific information regarding the analytical results. The data validation process is intended to evaluate the data on a technical basis. The data package also was subjected to an internal laboratory quality review prior to submission to Earth Tech for data validation.

During the validation process, laboratory-qualified and unqualified data are verified against all available supporting documentation. Based on this evaluation, qualifier codes may be added, deleted or modified by the data user. Final results are therefore, either qualified or unqualified. Validator-qualified results are annotated with the following codes in accordance with the Functional Guidelines:

- U The constituent was analyzed for, but was not detected above the level of the associated analytical reporting limit. The associated value is either the sample quantitation limit or the sample detection limit.

- J The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample.
- UJ The analyte was not detected above the reported sample quantitation limit. However, the reported quantitation limit is approximate and may or may not represent the actual limit of quantitation necessary to accurately and precisely measure the analyte in the sample.
- R The sample results are rejected due to serious deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte cannot be verified.

Details of the semivolatile data validation findings and conclusions are provided in the following sections of this report:

1. Holding Times
2. GC/MS Tuning
3. Calibration
 - A. IC
 - B. CC
4. Blanks
5. System Monitoring Compound Recovery
6. MS/MSD
7. Internal Standards Performance
8. Compound Identification
9. Constituent Quantitation and Reported Detection Limits
10. System Performance
11. Documentation
12. Overall Assessment

1. HOLDING TIMES

All samples were initially extracted within the seven-day technical holding time and the five-day VTSR method holding time. The cooler temperature upon receipt at the laboratory was within the recommended temperature of 4°C +/- 2°C.

2. GC/MS TUNING

The samples were analyzed on a single GC/MS system, identified as MSSV2. Two decafluorotriphenylphosphine (DFTPP) tunes were run representing the shift in which the standards and samples were analyzed. The DFTPP tunes are acceptable.

3. CALIBRATION

A. Initial Calibration

One IC dated 3/19/04 was analyzed in support of the semivolatile sample analyses. Documentation of the IC was present in the data package, and the Relative Response Factor (RRF), as well as percent % RSD values were accurately reported for all target compounds. The criteria employed for technical data review purposes are different than those used in the method. The laboratory must meet a minimum RRF of 0.01; however, for data review purposes, a RRF criterion of "greater than or equal to 0.05" is applied to all semi-volatile compounds. The RRF's and the average RRF were within the acceptance criteria specified in the method for all reported analytes. The %RSD's were within the acceptance criteria specified in the method for all target analytes with the exception of Di-n-butylphthalate (32.0%), Di-n-octylphthalate (30.3%) and Diethylphthalate (41.8%). The lowest point of the calibration curve was dropped for Di-n-butylphthalate, Di-n-octylphthalate, and Diethylphthalate and the %RSD were recalculated. The recalculated %RSD's were within the acceptance criteria of less than 30%. Di-n-octylphthalate and Diethylphthalate were not detected in the associated samples therefore data qualification was not required. The detected results for Di-n-butylphthalate were mitigated do to the presence of Di-n-butylphthalate in the associated method blank.

B. Continuing Calibration

One CC dated 3/26/04 was analyzed in support of the semivolatile sample analyses reported in the data submissions. The RRF's for the CC dated 3/26/04 were within the acceptance criteria. The percent difference (%D) between the average RRF's and the CC Response Factors for the CC dated 3/26/04 were within the acceptance criteria with the exception the %D for Di-n-butylphthalate and Di-n-octylphthalate. As per the National Functional Guidelines, if the %D exceeds the acceptance criteria qualify detected results for that analyte with "J" and non-detected results for that analyte with "UJ".

4. BLANKS

One laboratory semivolatile method blank and field blank were analyzed with this SDG. The results are summarized below.

Method Blank (0322SBLK)

Di-n-butylphthalate was detected at a concentration of 0.604 ppb in method blank 0322SBLK.

Field Blank (SKGWFB1009)

There were no analytes detected above the MDL in the field blank collected on 3/18/04.

5. SYSTEM MONITORING COMPOUND RECOVERY

All reported semivolatile system monitoring compounds were recovered within acceptable control limits.

6. MATRIX SPIKE/MATRIX SPIKE DUPLICATE (MS/MSD)

Sample SKGWD611009 was submitted for MS/MSD analysis. The MS/MSD percent recoveries were within the acceptance criteria with the exception of the 4-Nitorphenol. The %RPD between the MS/MSD are within the acceptance criteria. As per the National Functional Guidelines, no action is taken on MS/MSD data alone.

7. INTERNAL STANDARDS PERFORMANCE

Internal standard areas and retention times were within acceptable limits for the reported semivolatile sample analyses.

8. COMPOUND IDENTIFICATION

All reported semivolatile constituents were correctly identified with supporting chromatograms present in the data package.

9. CONSTITUENT QUANTITATION AND REPORTED DETECTION LIMITS

Constituent quantitations were correctly calculated and reported for semivolatile constituents.

10. SYSTEM PERFORMANCE

The analytical system appears to have been working well at the time of these analyses, based on the evaluation of the raw data submitted for review.

11. DOCUMENTATION

The documentation appeared accurate and in order with the exception of the following: GCAL inadvertently left the Date Extracted off of the CLP Form 1's therefore the data validator inserted the Date Extracted on the CLP Form 1's. GCAL also inadvertently left the "B" qualifier off of the CLP Form 1's for the compound Di-n-butylphthalate therefore the data validator inserted a "B" qualifier in the "Q" column of the CLP Form 1's.

The "B" qualifier indicates that the analyte was detected in the associated method blank. On pages 210/212/213/214 sample SKGW07R1009 was not listed but 2 field blanks were listed while only one field blank submitted for analysis. GCAL corrected the mistake and reissued corrected pages.

12. OVERALL ASSESSMENT

The results are acceptable with the validator-added qualifiers.

**DATA VALIDATION SUMMARY – SAMPLE DELIVERY GROUP 204031909
VOLATILE ORGANIC**

Validation of the GC/MS volatile organics data, as prepared by Gulf Coast Analytical Laboratories (GCAL) for the samples collected from the Skinner Landfill site in March 2004, was conducted by Earth Tech using the National Functional Guidelines for Organic Data Review, (US EPA, October, 1999), as appropriate. The results were reported by GCAL under SDG 204031909.

| GCAL # | Sample Description |
|-------------|--------------------|
| 20403190901 | SKGW581009 |
| 20403190902 | SKGW58DUP1009 |
| 20403190903 | SKGW631009 |
| 20403190904 | SKGW611009 |
| 20403190905 | SKGW61MS1009 |
| 20403190906 | SKGW61MSD1009 |
| 20403190908 | SKGWFB1009 |
| 20403190909 | SKTB1009 |
| 20403220801 | SKGW06R1009 |
| 20403220802 | SKGW07R1009 |
| 20403220803 | SKGW591009 |
| 20403220804 | SKGW601009 |
| 20403220805 | SKGW62A1009 |
| 20403220806 | SKGW641009 |
| 20403220807 | SKGW651009 |
| 20403220814 | TRIP BLANK |

INTRODUCTION

Analyses were performed according to CLP-Organic Analysis Low Concentration OLC02.0 SOW. Results of the sample analyses are reported by the laboratory as either qualified or unqualified. Unqualified results mean that the reported values may be used without reservation. Various qualifier codes are used by the laboratory to denote specific information regarding the analytical results. The data validation process is intended to evaluate the data on a technical basis. The data package also was subjected to an internal laboratory quality review prior to submission to Earth Tech for data validation.

During the validation process, laboratory-qualified and unqualified data are verified against all available supporting documentation. Based on this evaluation, qualifier codes may be added, deleted or modified by the data user. Final results are therefore, either qualified or unqualified. Validator-qualified results are annotated with the following codes in accordance with the Functional Guidelines:

- U The constituent was analyzed for, but was not detected above the level of the associated analytical reporting limit. The associated value is either the sample quantitation limit or the sample detection limit.
- J The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample.
- UJ The analyte was not detected above the reported sample quantitation limit.
- However, the reported quantitation limit is approximate and may or may not represent the actual limit of quantitation necessary to accurately and precisely measure the analyte in the sample.
- R The sample results are rejected due to serious deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte cannot be verified.

The volatiles data validation findings and conclusions are provided in the following sections of this report:

1. Holding Times
2. GC/MS Tuning
3. Calibration
 - A. IC
 - B. CC
4. Blanks
5. System Monitoring Compound Recovery
6. MS/MSD
7. Laboratory Control Sample
8. Internal Standards Performance
9. Compound Identification
10. Constituent Quantitation and Reported Detection Limits
11. System Performance

12. Documentation
13. Overall Assessment

1. HOLDING TIMES

All samples for Volatile Organic Compounds (VOC) analyses were analyzed within the 14-day technical holding time and the 10-day VTSR method holding time. The cooler temperature upon receipt at the laboratory was within the recommended temperature of 4°C +/- 2°C.

2. GC/MS TUNING

The samples were analyzed on two GC/MS system, identified as MSV0 and MSV2. Two bromofluorobenzene (BFB) tunes were run. The BFB tunes are acceptable.

3. CALIBRATION

A. Initial Calibration

Two IC's dated 3/21/04 and 3/23/04 were analyzed in support of the volatile sample analyses reported in the data submissions. The IC dated 3/21/04 was analyzed on instrument MSV0 and the IC dated 3/23/04 was analyzed on instrument MSV2. Documentation of the IC standards is present in the data package, and RRF's as well as %RSD values were accurately reported. The criteria employed for technical data review purposes are different than those used in the method. The laboratory must meet a minimum RRF of 0.01; however, for data review purposes, a RRF criterion of "greater than or equal to 0.05" is applied to all volatile compounds. The %RSD's were within the acceptance criteria specified in the method for all target analytes.

The RRF's and the average RRF for the IC dated 3/21/04 were within the acceptance criteria specified in the method for all target analytes with the exception of Acetone and 2-Butnaone. The RRF for the 1.0 ppb standard was below the acceptance criteria. The data validator dropped the 1.0 ppb standard for 1,2-Dibromo-3-chloropropane from the calibration curve and re-calculated the average RRF and %RSD. 1,2-Dibromo-3-chloropropane was not detected in any of the associated samples therefore data qualification was not required.

The RRF's and the average RRF for the IC dated 3/23/04 were within the acceptance criteria specified in the method for all target analytes with the exception of Acetone.

As per the National Functional Guidelines, if any initial calibration RRF is less than 0.05, qualify positive results that have acceptable mass spectral identification with "J", using professional judgement, and non-detected analytes as unusable (R).

It should be noted that the laboratory did meet the minimum RRF of 0.01 for all target compounds.

B. Continuing Calibration

Two CC dated 3/21/04 and 3/23/04 were analyzed on instrument MSV2 in support of the volatile sample analyses reported in the data submissions.

CC dated 3/21/04

The percent difference (%D) between the average RRF's and the CC RF's were within the acceptance criteria for all target analytes. The CC RRF's were within the acceptance criteria specified in the method for all target analytes with the exception of Acetone and 2-Butnaone. The Acetone and 2-Butnaone results were previously qualified under section 3A above.

CC dated 3/23/04

The percent difference (%D) between the average RRF's and the CC RF's were within the acceptance criteria for all target analytes. The CC RRF's were within the acceptance criteria specified in the method for all target analytes with the exception of Acetone. The Acetone results were previously qualified under section 3A above.

4. BLANKS

Two laboratory volatile method blanks, storage blank, two Trip Blanks, and a Field Blank were analyzed with this SDG. The results are summarized below.

MB153850

1,2-Dichlorobenzene and 1,3-Dichlorobenzene were detected at concentrations of 0.20 ppb and 0.12 ppb respectively in the method blank analyzed on 3/21/04.

MB153988

Acetone, 1,2-Dichlorobenzene and 1,3-Dichlorobenzene were detected at concentrations of 1.5 ppb, 0.21 ppb and 0.16 ppb respectively in the method blank analyzed on 3/23/04.

Storage Blank (CLP Storage Blank)

Methylene chloride, 1,4-Dichlorobenzene, and 1,2-Dichlorobenzene were detected at concentrations of 0.15 ppb, 0.096 ppb and 0.11 ppb respectively in the storage blank analyzed on 3/23/04.

Trip Blank (SKTB1009)

Methylene chloride was detected at a concentration of 0.23 ppb in the Trip Blank submitted for the sampling event that occurred on 3/18/04. The Methylene chloride detected in the trip blank was mitigated by the presence of Methylene chloride in the associated storage blank.

Trip Blank

Acetone and Methylene chloride were detected at concentrations 3.1 ppb and 0.24 ppb respectively in the Trip Blank submitted for the sampling event that occurred on 3/16-17/04. The Methylene chloride detected in the trip blank was mitigated by the presence of Methylene chloride in the associated storage blank.

Field Blank (SKGWFB1009)

Methylene chloride was detected at a concentration of 0.52 ppb in the Field Blank collected on 3/18/04. The Methylene chloride detected in the trip blank was mitigated by the presence of Methylene chloride in the associated storage blank.

5. SYSTEM MONITORING COMPOUND RECOVERY

All reported volatile system monitoring compounds were recovered within acceptable control limits for all samples.

6. MATRIX SPIKE/MATRIX SPIKE DUPLICATE

Sample SKGWD611009 was submitted for MS/MSD analysis. The MS/MSD percent recoveries (60%-140%) and %RPD (<40%) between the MS/MSD were within the acceptance criteria with the exception of the following:

| Compound | MS | MSD | RPD |
|----------------------|------------|------------|-----------|
| 1,1-Dichloroethene | 142 | 150 | 5 |
| 2-Hexanone | 58 | 87 | 40 |
| 4-methyl-2-pentanone | 67 | 102 | 41 |
| Bromomethane | 151 | 156 | 3 |
| Carbon disulfide | 145 | 145 | 0 |
| Chloroethane | 132 | 145 | 9 |
| Chloromethane | 135 | 146 | 8 |
| Vinyl chloride | 135 | 143 | 6 |

The results that are bolded exceeded the acceptance criteria. As per the National Functional Guidelines, no action is taken on MS/MSD data alone.

7. LABORATORY CONTROL SAMPLE

A LCS was analyzed in conjunction with this SDG. Recoveries were within the control limit for all constituents.

8. INTERNAL STANDARDS PERFORMANCE

Internal Standard (IS) areas and retention times were within acceptable limits for the reported volatile sample analyses.

9. COMPOUND IDENTIFICATION

All reported VOCs were correctly identified with supporting chromatograms present in the data package.

10. CONSTITUENT QUANTITATION AND REPORTED DETECTION LIMITS

Constituent quantitations were correctly calculated and reported for VOCs with the exception of Ethylbenzene. The Ethylbenzene standard and detected results were originally quantitated using the incorrect quantitation ion (GCAL used 106 instead of 91). GCAL corrected the mistake and re-submitted the corrected pages that were affected in the laboratory report. The overall effect had no impact in the final result for Ethylbenzene.

11. SYSTEM PERFORMANCE

The analytical system appears to have been working well at the time of these analyses, based on the evaluation of the raw data.

12. DOCUMENTATION

The documentation appeared accurate and in order with the exception of the following: GCAL inadvertently transposed the area counts and retention times for IS Chlorobenzene-d5 and 1,4-Dichlorobenzene-d4. The data validator corrected the mistake by drawing arrows to indicate the correct area counts and retention times for IS Chlorobenzene-d5 and 1,4-Dichlorobenzene-d4.

13. OVERALL ASSESSMENT

The results are acceptable with the validator-added qualifiers.

**DATA VALIDATION SUMMARY - SAMPLE DELIVERY GROUP 204031909
PESTICIDES**

Validation of the Gas Chromatography (GC) pesticides data, as prepared by Gulf Coast Analytical Laboratories (GCAL) for the samples collected from the Skinner Landfill site in March 2004, was conducted by Earth Tech using the National Functional Guidelines for Organic Data Review, (US EPA, October, 1999), as appropriate. The results were reported by GCAL under SDG 204031909.

| GCAL # | Sample Description |
|-------------|--------------------|
| 20403190901 | SKGW581009 |
| 20403190902 | SKGW58DUP1009 |
| 20403190903 | SKGW631009 |
| 20403190904 | SKGW611009 |
| 20403190905 | SKGW61MS1009 |
| 20403190906 | SKGW61MSD1009 |
| 20403190908 | SKGWFB1009 |
| 20403220801 | SKGW06R1009 |
| 20403220802 | SKGW07R1009 |
| 20403220803 | SKGW591009 |
| 20403220804 | SKGW601009 |
| 20403220805 | SKGW62A1009 |
| 20403220806 | SKGW641009 |

INTRODUCTION

Analyses were performed according to CLP-Organic Analysis Multi-Media, Multi-Concentration OLM04.2 SOW. Results of the sample analyses are reported by the laboratory as either qualified or unqualified. Unqualified results mean that the reported values may be used without reservation. Various qualifier codes are used by the laboratory to denote specific information regarding the analytical results.

The data validation process is intended to evaluate the data on a technical basis. The data package also was subjected to an internal laboratory quality review prior to submission to Earth Tech for data validation.

During the validation process, laboratory-qualified and unqualified data are verified against all available supporting documentation. Based on this evaluation, qualifier codes may be added, deleted or modified by the data user. Final results are therefore, either qualified or unqualified. Validator-qualified results are annotated with the following codes in accordance with the Functional Guidelines:

- U The constituent was analyzed for, but was not detected above the level of the associated analytical reporting limit. The associated value is either the sample quantitation limit or the sample detection limit.

- J The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample.
- UJ The analyte was not detected above the reported sample quantitation limit. However, the reported quantitation limit is approximate and may or may not represent the actual limit of quantitation necessary to accurately and precisely measure the analyte in the sample.
- R The sample results are rejected due to serious deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte cannot be verified.

Details of the pesticide data validation findings and conclusions are provided in the following sections of this report:

1. Holding Times
2. Gas Chromatograph/Electronic Capture Detector (GC/ECD) Instrument Performance Check
3. IC
4. Calibration Verification
5. Blanks
6. Surrogate Spikes
7. Matrix Spike/Matrix Spike Duplicate (MS/MSD)
8. Pesticide Cleanup Checks
9. Target Compound Identification
10. Constituent Quantitation and Reported Detection Limits
11. Documentation
12. Overall Assessment

1. HOLDING TIMES

All samples were extracted within the seven-day technical holding time and the five-day VTSR method holding time. The cooler temperature upon receipt at the laboratory was within the recommended temperature of 4°C +/- 2°C.

2. GC/ECD INSTRUMENT PERFORMANCE CHECK

The Performance Evaluation Mixture (PEM) was analyzed at the correct frequency. Absolute retention times were within limits.

The percent resolution between adjacent peaks was within QC limits for the Pesticide Analyte Resolution Check. The percent resolution between adjacent peaks is within QC limits for the Performance Evaluation Mixtures (PEM). The percent breakdown for both 4,4-DDT and Endrin in each PEM was less than 20.0% for both GC columns. The combined percent breakdown for 4,4-DDT and Endrin in each PEM was less than 30.0% for both GC columns.

3. INITIAL CALIBRATION

Individual standard mixtures A and B were analyzed at the correct frequencies and concentrations. The percent resolution criterion was met for Individual standard mixtures A and B.

The Percent Relative Standard Deviation (%RSD) of the calibration factors for each of the single component pesticides was less than 20%.

The multi-component target compounds were analyzed separately on both columns at a single concentration level. Retention times were determined from a minimum of three peaks.

4. CALIBRATION VERIFICATION

Absolute retention times were within appropriate time retention windows

5. BLANKS

One laboratory method blank and field blank were analyzed with this SDG. The results are summarized below.

Method Blank 154072

No constituents were detected above the laboratory-reporting limit. This blank corresponds to all samples extracted on 3/23/04

Field Blank (SKGWFB1009)

No constituents were detected above the laboratory-reporting limit in the field blank collected on 3/18/04.

6. SURROGATE SPIKES

Decachlorobiphenyl (DCB) and tetrachloro-m-xylene (TCX) surrogate spike recoveries were within the acceptance criteria for all samples.

7. MATRIX SPIKE/MATRIX SPIKE DUPLICATE

Sample SKGWD611009 was submitted for MS/MSD analysis. The MS/MSD percent recoveries were within the acceptance criteria. The %RPD between the MS/MSD are within the acceptance criteria.

8. PESTICIDE CLEANUP CHECKS

Recoveries of all pesticides and surrogates were within 80-120% for the lot of Florisil cartridges utilized for pesticide cleanup with the exception of 4,4'-DDT (130%). There were no target compounds detected in the associated samples therefore no action was taken.

9. TARGET COMPOUND IDENTIFICATION

All reported pesticide data were correctly identified with supporting chromatograms present in the data package.

10. CONSTITUENT QUANTITATION AND REPORTED DETECTION LIMITS

Constituent quantitations were correctly calculated and reported for pesticide constituents.

11. DOCUMENTATION

The documentation appeared accurate and in order with the exception of the following: GCAL inadvertently left the Date Extracted off of the CLP Form 1's therefore the data validator inserted the Date Extracted on the CLP Form 1's.

12. OVERALL ASSESSMENT

The results are acceptable with the validator-added qualifiers.

REFERENCES

US EPA, 1994. *National Functional Guidelines for Inorganic Data Review.*

US EPA, 1999. *National Functional Guidelines for Organic Data Review.*

ANALYTICAL RESULTS

PERFORMED BY

GULF COAST ANALYTICAL LABORATORIES, INC.

Report Date 04/08/2004

GCAL Report 204031909

ADDENDUM

Deliver To Earth Tech
200 Vine Street
Wilder, KY 41076
859-442-2300

Attn Pat Higgins

Customer Earth Tech

Project Skinner Landfill

000001

CASE NARRATIVE

Client: Earth Tech **Report:** 204031909

Gulf Coast Analytical Laboratories received and analyzed the sample(s) listed on the sample cross-reference page of this report. Receipt of the sample(s) is documented by the attached chain of custody. This applies only to the sample(s) listed in this report. No sample integrity or quality control exceptions were identified unless noted below.

The ILM04.1 – CLP analysis is resubmitted as an addendum to include an expanded list of compounds at the request of the client. The Addendum includes an updated case narrative.

VOLATILES MASS SPECTROMETRY

In analytical batch 271073, the MS/MSD exhibited sporadic recovery and RPD failures. The LCS/LCSD recoveries were acceptable.

In analytical batch 271108, no MS/MSD was analyzed due to insufficient sample volume. All LCS/LCSD recoveries were acceptable.

SEMI-VOLATILES MASS SPECTROMETRY

The MS/MSD recoveries for 4-Nitrophenol were above the upper control limit. All other batch QC was acceptable.

SEMI-VOLATILE GAS CHROMATOGRAPHY

In the Pesticide analysis, the recovery for DDT was above control limits in the Florisil check, however DDT was not detected in the associated samples.

METALS

In the ILM04.1 - CLP Metals analysis for prep batch 271122, the MS recovery was outside the control limits for Selenium. The LCS recovery was within control limits. This indicates the analysis is in control and the sample is in control and the sample is affected by matrix interference. A post-digestion spike was performed on the QC sample for this batch with a recovery of 49%. Aluminum, Copper, Lead, and Zinc is flagged as estimated for samples associated with prep batch 271122 due to the fact that the percent difference between the original sample result and the serial dilution result is greater than 10. A *chemical or physical interference is suspected. The Sample/Duplicate RPD for Vanadium for prep batch 271122 is not applicable because the sample and/or duplicate concentration is less than five times the reporting limit.*

In the ILM04.1 - CLP Metals analysis for prep batch 271124, the MS recovery was outside the control limits for Selenium. The LCS recovery was within control limits. This

000002
RESUBMITTED

indicates the analysis is in control and the sample is in control and the sample is affected by matrix interference. A post-digestion spike was performed on the QC sample for this batch with a recovery of 0%. Iron and Manganese is flagged as estimated for samples associated with prep batch 271124 due to the fact that the percent difference between the original sample result and the serial dilution result is greater than 10. A chemical or physical interference is suspected. The Sample/Duplicate RPD for Vanadium for prep batch 271124 is not applicable because the sample and/or duplicate concentration is less than five times the reporting limit.

The Sample/Duplicate RPD for Cyanide, Total for prep batch 271113 is not applicable because the sample and/or duplicate concentration is less than five times the reporting limit.

000002A

ADDITION

U.S. EPA - CLP
1
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKGW581009

Lab Name: PROJ AAH GCAL Contract: _____
 Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: _____
 Matrix: (soil / water) Water Lab Sample ID: 20403190901
 Level: (low / med) _____ Date Received: 03/19/04
 % Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|--------------|----|
| 7429-90-5 | Aluminum | 12000 | | E | P |
| 7440-36-0 | Antimony | 5.7 | B | | P |
| 7440-38-2 | Arsenic | 11.5 | | | P |
| 7440-39-3 | Barium | 284 | | | P |
| 7440-41-7 | Beryllium | 1.0 | B | | P |
| 7440-43-9 | Cadmium | 1.5 | B | | P |
| 7440-70-2 | Calcium | 214000 | | | P |
| 7440-47-3 | Chromium | 28.2 | | | P |
| 7440-48-4 | Cobalt | 13.4 | B | | P |
| 7440-50-8 | Copper | 45.7 | | E | P |
| 7439-89-6 | Iron | 32700 | | | P |
| 7439-92-1 | Lead | 19.5 | | E | P |
| 7439-95-4 | Magnesium | 56000 | | | P |
| 7439-96-5 | Manganese | 1300 | | | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 32.1 | B | | P |
| 7440-09-7 | Potassium | 7640 | | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-23-5 | Sodium | 33500 | | | P |
| 7440-28-0 | Thallium | 4.1 | B | | P |
| 7440-62-2 | Vanadium | 23.2 | B | | P |
| 7440-66-6 | Zinc | 81.0 | | E | P |
| 57-12-5 | Cyanide | 0.5 | U | | AS |

9/30/05
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Color Before: LT.BROWN Clarity Before: CLEAR Texture: _____
 Color After: LT.BROWN Clarity After: CLEAR Artifacts: _____
 Comments: _____

U.S. EPA - CLP
1
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKGW58DUP1009

Lab Name: PROJ AAH GCAL Contract: _____
 Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: _____
 Matrix: (soil / water) Water Lab Sample ID: 20403190902
 Level: (low / med) _____ Date Received: 03/19/04
 % Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|--------------|----|
| 7429-90-5 | Aluminum | 17100 | | P | P |
| 7440-36-0 | Antimony | 3.7 | B | | P |
| 7440-38-2 | Arsenic | 15.1 | | | P |
| 7440-39-3 | Barium | 326 | | | P |
| 7440-41-7 | Beryllium | 1.3 | B | | P |
| 7440-43-9 | Cadmium | 1.9 | B | | P |
| 7440-70-2 | Calcium | 256000 | | | P |
| 7440-47-3 | Chromium | 41.8 | | | P |
| 7440-48-4 | Cobalt | 19.5 | B | | P |
| 7440-50-8 | Copper | 55.9 | | P | P |
| 7439-89-6 | Iron | 48400 | | | P |
| 7439-92-1 | Lead | 27.2 | | P | P |
| 7439-95-4 | Magnesium | 63200 | | | P |
| 7439-96-5 | Manganese | 1500 | | | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 46.4 | | | P |
| 7440-09-7 | Potassium | 8480 | | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-23-5 | Sodium | 34600 | | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-62-2 | Vanadium | 32.9 | B | | P |
| 7440-66-6 | Zinc | 125 | | P | P |
| 57-12-5 | Cyanide | 0.5 | U | | AS |

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Color Before: LT.BROWN Clarity Before: CLEAR Texture: _____
 Color After: LT.BROWN Clarity After: CLEAR Artifacts: _____
 Comments: _____

U.S. EPA - CLP
1
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKGW631009

Lab Name: PROJ AAH GCAL

Contract: _____

Lab Code: LA024

Case No.: _____

SAS No.: _____

SDG No.: _____

Matrix: (soil / water) Water

Lab Sample ID: 20403190903

Level: (low / med) _____

Date Received: 03/19/04

% Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|-------------------------------------|----|
| 7429-90-5 | Aluminum | 26600 | | <input checked="" type="checkbox"/> | P |
| 7440-36-0 | Antimony | 5.7 | B | | P |
| 7440-38-2 | Arsenic | 17.1 | | | P |
| 7440-39-3 | Barium | 186 | B | | P |
| 7440-41-7 | Beryllium | 2.1 | B | | P |
| 7440-43-9 | Cadmium | 2.5 | B | | P |
| 7440-70-2 | Calcium | 465000 | | | P |
| 7440-47-3 | Chromium | 38.2 | | | P |
| 7440-48-4 | Cobalt | 28.3 | B | | P |
| 7440-50-8 | Copper | 69.2 | | <input checked="" type="checkbox"/> | P |
| 7439-89-6 | Iron | 63200 | | | P |
| 7439-92-1 | Lead | 41.0 | | <input checked="" type="checkbox"/> | P |
| 7439-95-4 | Magnesium | 111000 | | | P |
| 7439-96-5 | Manganese | 2570 | | | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 58.1 | | | P |
| 7440-09-7 | Potassium | 9320 | | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-23-5 | Sodium | 45000 | | | P |
| 7440-28-0 | Thallium | 8.5 | B | | P |
| 7440-62-2 | Vanadium | 43.0 | B | | P |
| 7440-66-6 | Zinc | 176 | | <input checked="" type="checkbox"/> | P |
| 57-12-5 | Cyanide | 0.5 | U | | AS |

Color Before: LT.BROWN

Clarity Before: CLEAR

Texture: _____

Color After: LT.BROWN

Clarity After: CLEAR

Artifacts: _____

Comments:

U.S. EPA - CLP
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INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKGW611009

Lab Name: PROJ AAH GCAL Contract: _____

Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: _____

Matrix: (soil / water) Water Lab Sample ID: 20403190904

Level: (low / med) _____ Date Received: 03/19/04

% Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|--------------|----|
| 7429-90-5 | Aluminum | 452 | | E | P |
| 7440-36-0 | Antimony | 4.8 | B | | P |
| 7440-38-2 | Arsenic | 2.9 | U | | P |
| 7440-39-3 | Barium | 44.1 | B | | P |
| 7440-41-7 | Beryllium | 0.2 | B | | P |
| 7440-43-9 | Cadmium | 0.3 | B | | P |
| 7440-70-2 | Calcium | 187000 | | | P |
| 7440-47-3 | Chromium | 1.9 | B | | P |
| 7440-48-4 | Cobalt | 1.7 | B | | P |
| 7440-50-8 | Copper | 22.2 | B | E | P |
| 7439-89-6 | Iron | 2430 | | | P |
| 7439-92-1 | Lead | 22.1 | | E | P |
| 7439-95-4 | Magnesium | 30000 | | | P |
| 7439-96-5 | Manganese | 527 | | | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 4.3 | B | | P |
| 7440-09-7 | Potassium | 6950 | | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-23-5 | Sodium | 27000 | | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-62-2 | Vanadium | 2.1 | B | | P |
| 7440-66-6 | Zinc | 7.3 | B | E | P |
| 57-12-5 | Cyanide | 0.5 | U | | AS |

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Color Before: COLORLESS

Clarity Before: CLEAR

Texture: _____

Color After: COLORLESS

Clarity After: CLEAR

Artifacts: _____

Comments:

U.S. EPA - CLP
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INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKGW61MS1009

Lab Name: PROJ AAH GCAL Contract: _____
 Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: _____
 Matrix: (soil / water) Water Lab Sample ID: 20403190905
 Level: (low / med) _____ Date Received: 03/19/04
 % Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|----------|----|
| 7429-90-5 | Aluminum | 2520 | | <i>E</i> | P |
| 7440-36-0 | Antimony | 116 | | | P |
| 7440-38-2 | Arsenic | 50.0 | | | P |
| 7440-39-3 | Barium | 2120 | | | P |
| 7440-41-7 | Beryllium | 53.8 | | | P |
| 7440-43-9 | Cadmium | 50.4 | | | P |
| 7440-70-2 | Calcium | 188000 | | | P |
| 7440-47-3 | Chromium | 209 | | | P |
| 7440-48-4 | Cobalt | 498 | | | P |
| 7440-50-8 | Copper | 279 | | <i>E</i> | P |
| 7439-89-6 | Iron | 3390 | | | P |
| 7439-92-1 | Lead | 42.2 | | <i>E</i> | P |
| 7439-95-4 | Magnesium | 29400 | | | P |
| 7439-96-5 | Manganese | 1060 | | | P |
| 7439-97-6 | Mercury | 5.1 | | | AV |
| 7440-02-0 | Nickel | 503 | | | P |
| 7440-09-7 | Potassium | 7060 | | | P |
| 7782-49-2 | Selenium | 4.5 | B | N | P |
| 7440-22-4 | Silver | 51.5 | | | P |
| 7440-23-5 | Sodium | 27200 | | | P |
| 7440-28-0 | Thallium | 48.3 | | | P |
| 7440-62-2 | Vanadium | 531 | | | P |
| 7440-66-6 | Zinc | 480 | | <i>E</i> | P |
| 57-12-5 | Cyanide | 86.1 | | | AS |

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Color Before: _____ Clarity Before: _____ Texture: _____
 Color After: _____ Clarity After: _____ Artifacts: _____
 Comments: _____

U.S. EPA - CLP
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INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKGW61DUP1009

Lab Name: PROJ AAH GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: _____
Matrix: (soil / water) Water Lab Sample ID: 20403190907
Level: (low / med) _____ Date Received: 03/19/04
% Solids: _____
Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|--------------|----|
| 7429-90-5 | Aluminum | 470 | | E | P |
| 7440-36-0 | Antimony | 5.5 | B | | P |
| 7440-38-2 | Arsenic | 3.7 | B | | P |
| 7440-39-3 | Barium | 45.4 | B | | P |
| 7440-41-7 | Beryllium | 0.2 | B | | P |
| 7440-43-9 | Cadmium | 0.3 | B | | P |
| 7440-70-2 | Calcium | 192000 | | | P |
| 7440-47-3 | Chromium | 1.6 | B | | P |
| 7440-48-4 | Cobalt | 1.6 | B | | P |
| 7440-50-8 | Copper | 18.6 | B | E | P |
| 7439-89-6 | Iron | 2490 | | | P |
| 7439-92-1 | Lead | 22.5 | | E | P |
| 7439-95-4 | Magnesium | 30300 | | | P |
| 7439-96-5 | Manganese | 530 | | | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 3.7 | B | | P |
| 7440-09-7 | Potassium | 7140 | | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-23-5 | Sodium | 27800 | | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-62-2 | Vanadium | 1.7 | B | | P |
| 7440-66-6 | Zinc | 7.0 | B | E | P |
| 57-12-5 | Cyanide | 0.8 | B | | AS |

Color Before: _____ Clarity Before: _____ Texture: _____
Color After: _____ Clarity After: _____ Artifacts: _____
Comments: _____

U.S. EPA - CLP
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INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKGWFB1009

Lab Name: PROJ AAH GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: _____
Matrix: (soil / water) Water Lab Sample ID: 20403190908
Level: (low / med) _____ Date Received: 03/19/04
% Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|--------------|----|
| 7429-90-5 | Aluminum | 25.8 | U | P | P |
| 7440-36-0 | Antimony | 3.7 | U | | P |
| 7440-38-2 | Arsenic | 2.9 | U | | P |
| 7440-39-3 | Barium | 0.4 | B | | P |
| 7440-41-7 | Beryllium | 0.1 | U | | P |
| 7440-43-9 | Cadmium | 0.2 | U | | P |
| 7440-70-2 | Calcium | 93.4 | B | | P |
| 7440-47-3 | Chromium | 0.8 | U | | P |
| 7440-48-4 | Cobalt | 0.4 | U | | P |
| 7440-50-8 | Copper | 1.2 | U | P | P |
| 7439-89-6 | Iron | 14.1 | U | | P |
| 7439-92-1 | Lead | 1.5 | U | P | P |
| 7439-95-4 | Magnesium | 72.5 | B | | P |
| 7439-96-5 | Manganese | 0.8 | B | | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 0.9 | B | | P |
| 7440-09-7 | Potassium | 42.1 | U | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-23-5 | Sodium | 139 | B | | P |
| 7440-28-0 | Thallium | 4.6 | B | | P |
| 7440-62-2 | Vanadium | 0.8 | U | | P |
| 7440-66-6 | Zinc | 3.2 | B | P | P |
| 57-12-5 | Cyanide | 0.5 | U | | AS |

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Color Before: COLORLESS Clarity Before: CLEAR Texture: _____
Color After: COLORLESS Clarity After: CLEAR Artifacts: _____
Comments: _____

U.S. EPA - CLP
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INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKGW581009(DISS)

Lab Name: PROJ AAH GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: _____
Matrix: (soil / water) Water Lab Sample ID: 20403190910
Level: (low / med) _____ Date Received: 03/19/04
% Solids: _____
Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum | 25.8 | U | | P |
| 7440-36-0 | Antimony | 3.7 | U | | P |
| 7440-38-2 | Arsenic | 3.1 | B | | P |
| 7440-39-3 | Barium | 156 | B | | P |
| 7440-41-7 | Beryllium | 0.1 | U | | P |
| 7440-43-9 | Cadmium | 0.2 | U | | P |
| 7440-70-2 | Calcium | 109000 | | | P |
| 7440-47-3 | Chromium | 1.5 | B | | P |
| 7440-48-4 | Cobalt | 1.3 | B | | P |
| 7440-50-8 | Copper | 2.9 | B | | P |
| 7439-89-6 | Iron | 209 | | | P |
| 7439-92-1 | Lead | 1.5 | U | | P |
| 7439-95-4 | Magnesium | 32500 | | | P |
| 7439-96-5 | Manganese | 549 | | E | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 2.6 | B | | P |
| 7440-09-7 | Potassium | 4550 | B | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-23-5 | Sodium | 32400 | | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-62-2 | Vanadium | 1.6 | B | | P |
| 7440-66-6 | Zinc | 0.6 | U | | P |

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Color Before: COLORLESS Clarity Before: CLEAR Texture: _____
Color After: COLORLESS Clarity After: CLEAR Artifacts: _____
Comments: _____

U.S. EPA - CLP
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INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKGW58DUP1009(DISS)

Lab Name: PROJ AAH GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: _____
Matrix: (soil / water) Water Lab Sample ID: 20403190911
Level: (low / med) _____ Date Received: 03/19/04
% Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|--------------|----|
| 7429-90-5 | Aluminum | 25.8 | U | | P |
| 7440-36-0 | Antimony | 3.7 | U | | P |
| 7440-38-2 | Arsenic | 2.9 | U | | P |
| 7440-39-3 | Barium | 164 | B | | P |
| 7440-41-7 | Beryllium | 0.1 | U | | P |
| 7440-43-9 | Cadmium | 0.2 | U | | P |
| 7440-70-2 | Calcium | 109000 | | | P |
| 7440-47-3 | Chromium | 1.3 | B | | P |
| 7440-48-4 | Cobalt | 1.3 | B | | P |
| 7440-50-8 | Copper | 1.2 | U | | P |
| 7439-89-6 | Iron | 201 | | P | P |
| 7439-92-1 | Lead | 1.5 | U | | P |
| 7439-95-4 | Magnesium | 34400 | | | P |
| 7439-96-5 | Manganese | 492 | | E | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 2.0 | B | | P |
| 7440-09-7 | Potassium | 5130 | | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-23-5 | Sodium | 35200 | | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-62-2 | Vanadium | 1.4 | B | | P |
| 7440-66-6 | Zinc | 0.6 | U | | P |

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Color Before: COLORLESS Clarity Before: CLEAR Texture: _____
Color After: COLORLESS Clarity After: CLEAR Artifacts: _____
Comments:

U.S. EPA - CLP
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INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKGW631009(DISS)

Lab Name: PROJ AAH GCAL Contract: _____
 Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: _____
 Matrix: (soil / water) Water Lab Sample ID: 20403190912
 Level: (low / med) _____ Date Received: 03/19/04
 % Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|--------------|----|
| 7423-90-5 | Aluminum | 25.8 | U | | P |
| 7440-36-0 | Antimony | 3.8 | B | | P |
| 7440-38-2 | Arsenic | 2.9 | U | | P |
| 7440-39-3 | Barium | 20.1 | B | | P |
| 7440-41-7 | Beryllium | 0.2 | B | | P |
| 7440-43-9 | Cadmium | 0.2 | U | | P |
| 7440-70-2 | Calcium | 295000 | | | P |
| 7440-47-3 | Chromium | 1.8 | B | | P |
| 7440-48-4 | Cobalt | 1.1 | B | | P |
| 7440-50-8 | Copper | 2.0 | B | | P |
| 7439-89-6 | Iron | 21.4 | B | E | P |
| 7439-92-1 | Lead | 1.5 | U | | P |
| 7439-95-4 | Magnesium | 67000 | | | P |
| 7439-96-5 | Manganese | 271 | | E | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 3.2 | B | | P |
| 7440-09-7 | Potassium | 5210 | | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-23-5 | Sodium | 46100 | | | P |
| 7440-28-0 | Thallium | 4.6 | B | | P |
| 7440-62-2 | Vanadium | 0.8 | U | | P |
| 7440-66-6 | Zinc | 0.6 | U | | P |

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Color Before: COLORLESS Clarity Before: CLEAR Texture: _____
 Color After: COLORLESS Clarity After: CLEAR Artifacts: _____
 Comments: _____

U.S. EPA - CLP
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INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKGW611009(DISS)

Lab Name: PROJ AAH GCAL

Contract: _____

Lab Code: LA024

Case No.: _____

SAS No.: _____

SDG No.: _____

Matrix: (soil / water) Water

Lab Sample ID: 20403190913

Level: (low / med) _____

Date Received: 03/19/04

% Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum | 25.8 | U | | P |
| 7440-36-0 | Antimony | 4.5 | B | | P |
| 7440-38-2 | Arsenic | 2.9 | U | | P |
| 7440-39-3 | Barium | 39.4 | B | | P |
| 7440-41-7 | Beryllium | 0.1 | B | | P |
| 7440-43-9 | Cadmium | 0.3 | B | | P |
| 7440-70-2 | Calcium | 191000 | | | P |
| 7440-47-3 | Chromium | 1.1 | B | | P |
| 7440-48-4 | Cobalt | 1.4 | B | | P |
| 7440-50-8 | Copper | 8.0 | B | | P |
| 7439-89-6 | Iron | 187 | | | P |
| 7439-92-1 | Lead | 1.5 | U | | P |
| 7439-95-4 | Magnesium | 29100 | | | P |
| 7439-96-5 | Manganese | 485 | | E | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 4.2 | B | | P |
| 7440-09-7 | Potassium | 6990 | | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-23-5 | Sodium | 27900 | | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-62-2 | Vanadium | 1.2 | B | | P |
| 7440-66-6 | Zinc | 0.6 | U | | P |

Color Before: COLORLESS

Clarity Before: CLEAR

Texture: _____

Color After: COLORLESS

Clarity After: CLEAR

Artifacts: _____

Comments: _____

U.S. EPA - CLP
1
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKGW61MS1009(DISS)

Lab Name: PROJ AAH GCAL

Contract: _____

Lab Code: LA024

Case No.: _____

SAS No.: _____

SDG No.: _____

Matrix: (soil / water) Water

Lab Sample ID: 20403190914

Level: (low / med) _____

Date Received: 03/19/04

% Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|--------------|----|
| 7429-90-5 | Aluminum | 2150 | | | P |
| 7440-36-0 | Antimony | 117 | | | P |
| 7440-38-2 | Arsenic | 48.4 | | | P |
| 7440-39-3 | Barium | 2110 | | | P |
| 7440-41-7 | Beryllium | 54.3 | | | P |
| 7440-43-9 | Cadmium | 50.1 | | | P |
| 7440-70-2 | Calcium | 184000 | | | P |
| 7440-47-3 | Chromium | 208 | | | P |
| 7440-48-4 | Cobalt | 489 | | | P |
| 7440-50-8 | Copper | 272 | | | P |
| 7439-89-6 | Iron | 1210 | | P | P |
| 7439-92-1 | Lead | 18.7 | | | P |
| 7439-95-4 | Magnesium | 28600 | | | P |
| 7439-96-5 | Manganese | 991 | | E | P |
| 7439-97-6 | Mercury | 4.1 | | | AV |
| 7440-02-0 | Nickel | 491 | | | P |
| 7440-09-7 | Potassium | 6730 | | | P |
| 7782-49-2 | Selenium | 7.3 | | N | P |
| 7440-22-4 | Silver | 52.9 | | | P |
| 7440-23-5 | Sodium | 27300 | | | P |
| 7440-28-0 | Thallium | 53.5 | | | P |
| 7440-62-2 | Vanadium | 532 | | | P |
| 7440-66-6 | Zinc | 467 | | | P |

J

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9/2/05
msk

Color Before: _____

Clarity Before: _____

Texture: _____

Color After: _____

Clarity After: _____

Artifacts: _____

Comments: _____

U.S. EPA - CLP
1
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKGW61DUP1009 (DISS)

Lab Name: PROJ AAH GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: _____
Matrix: (soil / water) Water Lab Sample ID: 20403190915
Level: (low / med) _____ Date Received: 03/19/04
% Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|--------------|----|
| 7429-90-5 | Aluminum | 25.8 | U | | P |
| 7440-36-0 | Antimony | 3.7 | U | | P |
| 7440-38-2 | Arsenic | 2.9 | U | | P |
| 7440-39-3 | Barium | 38.9 | B | | P |
| 7440-41-7 | Beryllium | 0.1 | B | | P |
| 7440-43-9 | Cadmium | 0.2 | U | | P |
| 7440-70-2 | Calcium | 186000 | | | P |
| 7440-47-3 | Chromium | 1.3 | B | | P |
| 7440-48-4 | Cobalt | 1.3 | B | | P |
| 7440-50-8 | Copper | 5.2 | B | | P |
| 7439-89-6 | Iron | 195 | | P | P |
| 7439-92-1 | Lead | 1.5 | U | | P |
| 7439-95-4 | Magnesium | 28400 | | | P |
| 7439-96-5 | Manganese | 477 | | E | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 3.6 | B | | P |
| 7440-09-7 | Potassium | 6860 | | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-23-5 | Sodium | 27200 | | | P |
| 7440-28-0 | Thallium | 5.6 | B | | P |
| 7440-62-2 | Vanadium | 1.8 | B | | P |
| 7440-66-6 | Zinc | 0.6 | U | | P |

9/30/05
m

Color Before: _____ Clarity Before: _____ Texture: _____
Color After: _____ Clarity After: _____ Artifacts: _____
Comments: _____

U.S. EPA - CLP
1
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKGWFB1009 (DISS)

Lab Name: PROJ AAH GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: _____
Matrix: (soil / water) Water Lab Sample ID: 20403190916
Level: (low / med) _____ Date Received: 03/19/04
% Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|--------------|----|
| 7429-90-5 | Aluminum | 25.8 | U | | P |
| 7440-36-0 | Antimony | 3.7 | B | | P |
| 7440-38-2 | Arsenic | 2.9 | U | | P |
| 7440-39-3 | Barium | 0.3 | U | | P |
| 7440-41-7 | Beryllium | 0.1 | U | | P |
| 7440-43-9 | Cadmium | 0.2 | U | | P |
| 7440-70-2 | Calcium | 34.9 | B | | P |
| 7440-47-3 | Chromium | 0.8 | U | | P |
| 7440-48-4 | Cobalt | 0.5 | B | | P |
| 7440-50-8 | Copper | 1.2 | U | | P |
| 7439-89-6 | Iron | 14.1 | U | E | P |
| 7439-92-1 | Lead | 1.5 | U | | P |
| 7439-95-4 | Magnesium | 47.3 | B | | P |
| 7439-96-5 | Manganese | 0.2 | U | E | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 0.7 | U | | P |
| 7440-09-7 | Potassium | 42.1 | U | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-23-5 | Sodium | 120 | B | | P |
| 7440-28-0 | Thallium | 7.5 | B | | P |
| 7440-62-2 | Vanadium | 0.8 | U | | P |
| 7440-66-6 | Zinc | 0.6 | U | | P |

4J

J

4J

9/30/05
m

Color Before: COLORLESS Clarity Before: CLEAR Texture: _____
Color After: COLORLESS Clarity After: CLEAR Artifacts: _____
Comments: _____

U.S. EPA - CLP
1
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKGW06R1009

Lab Name: PROJ AAH GCAL

Contract: _____

Lab Code: LA024

Case No.: _____

SAS No.: _____

SDG No.: _____

Matrix: (soil / water) Water

Lab Sample ID: 20403220801

Level: (low / med) _____

Date Received: 03/19/04

% Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|--------------|----|
| 7429-90-5 | Aluminum | 9900 | | P | P |
| 7440-36-0 | Antimony | 5.5 | B | | P |
| 7440-38-2 | Arsenic | 12.4 | | | P |
| 7440-39-3 | Barium | 440 | | | P |
| 7440-41-7 | Beryllium | 1.1 | B | | P |
| 7440-43-9 | Cadmium | 1.0 | B | | P |
| 7440-70-2 | Calcium | 309000 | | | P |
| 7440-47-3 | Chromium | 16.9 | | | P |
| 7440-48-4 | Cobalt | 12.3 | B | | P |
| 7440-50-8 | Copper | 39.3 | | P | P |
| 7439-89-6 | Iron | 25300 | | | P |
| 7439-92-1 | Lead | 23.9 | | P | P |
| 7439-95-4 | Magnesium | 83600 | | | P |
| 7439-96-5 | Manganese | 988 | | | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 23.4 | B | | P |
| 7440-09-7 | Potassium | 3970 | B | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-23-5 | Sodium | 21900 | | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-62-2 | Vanadium | 22.2 | B | | P |
| 7440-66-6 | Zinc | 72.9 | | P | P |
| 57-12-5 | Cyanide | 1.0 | B | | AS |

9/30/05
M 22

Color Before: LT. YELLOW

Clarity Before: CLEAR

Texture: _____

Color After: LT. YELLOW

Clarity After: CLEAR

Artifacts: _____

Comments:

U.S. EPA - CLP
1
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKGW07R1009

Lab Name: PROJ AAH GCAL Contract: _____
 Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: _____
 Matrix: (soil / water) Water Lab Sample ID: 20403220802
 Level: (low / med) _____ Date Received: 03/19/04
 % Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|--------------|----|
| 7429-90-5 | Aluminum | 7810 | | F | P |
| 7440-36-0 | Antimony | 6.6 | B | | P |
| 7440-38-2 | Arsenic | 6.9 | B | | P |
| 7440-39-3 | Barium | 484 | | | P |
| 7440-41-7 | Beryllium | 0.8 | B | | P |
| 7440-43-9 | Cadmium | 0.9 | B | | P |
| 7440-70-2 | Calcium | 281000 | | | P |
| 7440-47-3 | Chromium | 12.9 | | | P |
| 7440-48-4 | Cobalt | 7.0 | B | | P |
| 7440-50-8 | Copper | 35.5 | | F | P |
| 7439-89-6 | Iron | 20200 | | | P |
| 7439-92-1 | Lead | 9.2 | | F | P |
| 7439-95-4 | Magnesium | 54000 | | | P |
| 7439-96-5 | Manganese | 1590 | | | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 17.8 | B | | P |
| 7440-09-7 | Potassium | 4510 | B | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-23-5 | Sodium | 31200 | | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-62-2 | Vanadium | 15.3 | B | | P |
| 7440-66-6 | Zinc | 51.2 | | F | P |
| 57-12-5 | Cyanide | 1.5 | B | | AS |

9/30/05
msk

Color Before: LT. YELLOW Clarity Before: CLEAR Texture: _____
 Color After: LT. YELLOW Clarity After: CLEAR Artifacts: _____
 Comments:

U.S. EPA - CLP
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INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKGW591009

Lab Name: PROJ AAH GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: _____
Matrix: (soil / water) Water Lab Sample ID: 20403220803
Level: (low / med) _____ Date Received: 03/19/04
% Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|--------------|----|
| 7429-90-5 | Aluminum | 816 | | E | P |
| 7440-36-0 | Antimony | 4.7 | B | | P |
| 7440-38-2 | Arsenic | 2.9 | U | | P |
| 7440-39-3 | Barium | 55.0 | B | | P |
| 7440-41-7 | Beryllium | 0.2 | B | | P |
| 7440-43-9 | Cadmium | 0.2 | U | | P |
| 7440-70-2 | Calcium | 243000 | | | P |
| 7440-47-3 | Chromium | 5.5 | B | | P |
| 7440-48-4 | Cobalt | 2.1 | B | | P |
| 7440-50-8 | Copper | 10.1 | B | E | P |
| 7439-89-6 | Iron | 3020 | | | P |
| 7439-92-1 | Lead | 1.5 | U | E | P |
| 7439-95-4 | Magnesium | 51500 | | | P |
| 7439-96-5 | Manganese | 224 | | | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 6.7 | B | | P |
| 7440-09-7 | Potassium | 32500 | | | P |
| 7732-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-23-5 | Sodium | 162000 | | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-62-2 | Vanadium | 2.3 | B | | P |
| 7440-66-6 | Zinc | 7.9 | B | E | P |
| 57-12-5 | Cyanide | 1.0 | B | | AS |

9/30/04
min

Color Before: DK.BROWN Clarity Before: CLEAR Texture: _____
Color After: DK.BROWN Clarity After: CLEAR Artifacts: _____
Comments:

U.S. EPA - CLP
1
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKGW601009

Lab Name: PROJ AAH GCAL

Contract: _____

Lab Code: LA024

Case No.: _____

SAS No.: _____

SDG No.: _____

Matrix: (soil / water) Water

Lab Sample ID: 20403220804

Level: (low / med) _____

Date Received: 03/19/04

% Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|--------------|----|
| 7429-90-5 | Aluminum | 32500 | | E | P |
| 7440-36-0 | Antimony | 9.7 | B | | P |
| 7440-38-2 | Arsenic | 17.0 | | | P |
| 7440-39-3 | Barium | 129 | B | | P |
| 7440-41-7 | Beryllium | 2.5 | B | | P |
| 7440-43-9 | Cadmium | 2.8 | B | | P |
| 7440-70-2 | Calcium | 492000 | | | P |
| 7440-47-3 | Chromium | 59.6 | | | P |
| 7440-48-4 | Cobalt | 36.1 | B | | P |
| 7440-50-8 | Copper | 54.5 | | E | P |
| 7439-89-6 | Iron | 74200 | | | P |
| 7439-92-1 | Lead | 40.4 | | E | P |
| 7439-95-4 | Magnesium | 112000 | | | P |
| 7439-96-5 | Manganese | 1410 | | | P |
| 7439-97-6 | Mercury | 0.1 | B | | AV |
| 7440-02-0 | Nickel | 67.3 | | | P |
| 7440-09-7 | Potassium | 11800 | | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-23-5 | Sodium | 44600 | | | P |
| 7440-28-0 | Thallium | 11.0 | | | P |
| 7440-62-2 | Vanadium | 51.2 | | | P |
| 7440-66-6 | Zinc | 180 | | E | P |

7/3/05
mk

Cclor Before: LT.BROWN

Clarity Before: CLEAR

Texture: _____

Cclor After: LT.BROWN

Clarity After: CLEAR

Artifacts: _____

Ccmmnts: _____

U.S. EPA - CLP
1
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKGW62A1009

Lab Name: PROJ AAH GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: _____
Matrix: (soil / water) Water Lab Sample ID: 20403220805
Level: (low / med) _____ Date Received: 03/19/04
% Solids: _____
Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|--------------|----|
| 7429-90-5 | Aluminum | 13200 | | E | P |
| 7440-36-0 | Antimony | 6.1 | B | | P |
| 7440-38-2 | Arsenic | 8.3 | B | | P |
| 7440-39-3 | Barium | 361 | | | P |
| 7440-41-7 | Beryllium | 1.1 | B | | P |
| 7440-43-9 | Cadmium | 1.6 | B | | P |
| 7440-70-2 | Calcium | 337000 | | | P |
| 7440-47-3 | Chromium | 29.6 | | | P |
| 7440-48-4 | Cobalt | 15.6 | B | | P |
| 7440-50-8 | Copper | 42.7 | | E | P |
| 7439-89-6 | Iron | 35000 | | | P |
| 7439-92-1 | Lead | 39.5 | | E | P |
| 7439-95-4 | Magnesium | 88000 | | | P |
| 7439-96-5 | Manganese | 1460 | | | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 35.4 | B | | P |
| 7440-09-7 | Potassium | 13900 | | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-23-5 | Sodium | 123000 | | | P |
| 7440-28-0 | Thallium | 6.9 | B | | P |
| 7440-62-2 | Vanadium | 23.0 | B | | P |
| 7440-66-6 | Zinc | 101 | | E | P |
| 57-12-5 | Cyanide | 1.0 | B | | AS |

Color Before: COLORLESS Clarity Before: CLEAR Texture: _____
Color After: COLORLESS Clarity After: CLEAR Artifacts: _____
Comments:

U.S. EPA - CLP
1
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKGW641009

Lab Name: PROJ AAH GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: _____
Matrix: (soil / water) Water Lab Sample ID: 20403220806
Level: (low / med) _____ Date Received: 03/19/04
% Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|----------|----|
| 7429-90-5 | Aluminum | 3080 | | <u>E</u> | P |
| 7440-36-0 | Antimony | 4.9 | B | | P |
| 7440-38-2 | Arsenic | 2.9 | U | | P |
| 7440-39-3 | Barium | 37.1 | B | | P |
| 7440-41-7 | Beryllium | 0.3 | B | | P |
| 7440-43-9 | Cadmium | 0.2 | U | | P |
| 7440-70-2 | Calcium | 213000 | | | P |
| 7440-47-3 | Chromium | 7.0 | B | | P |
| 7440-48-4 | Cobalt | 5.4 | B | | P |
| 7440-50-8 | Copper | 11.3 | B | <u>E</u> | P |
| 7439-89-6 | Iron | 7520 | | | P |
| 7439-92-1 | Lead | 1.5 | U | <u>E</u> | P |
| 7439-95-4 | Magnesium | 66000 | | | P |
| 7439-96-5 | Manganese | 1650 | | | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 16.4 | B | | P |
| 7440-09-7 | Potassium | 15000 | | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-23-5 | Sodium | 59800 | | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-62-2 | Vanadium | 5.3 | B | | P |
| 7440-66-6 | Zinc | 13.6 | B | <u>E</u> | P |
| 57-12-5 | Cyanide | 1.3 | B | | AS |

4J

J

9/30/04
mk

Color Before: COLORLESS Clarity Before: CLEAR Texture: _____
Color After: COLORLESS Clarity After: CLEAR Artifacts: _____
Comments: _____

U.S. EPA - CLP
1
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKGW06R1009 (DISS)

Lab Name: PROJ AAH GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: _____
Matrix: (soil / water) Water Lab Sample ID: 20403220808
Level: (low / med) _____ Date Received: 03/19/04
% Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum | 25.8 | U | | P |
| 7440-36-0 | Antimony | 3.7 | U | | P |
| 7440-38-2 | Arsenic | 2.9 | U | | P |
| 7440-39-3 | Barium | 266 | | | P |
| 7440-41-7 | Beryllium | 0.1 | U | | P |
| 7440-43-9 | Cadmium | 0.2 | U | | P |
| 7440-70-2 | Calcium | 189000 | | | P |
| 7440-47-3 | Chromium | 1.2 | B | | P |
| 7440-48-4 | Cobalt | 0.4 | U | | P |
| 7440-50-8 | Copper | 1.2 | U | | P |
| 7439-89-6 | Iron | 22.0 | B | | P |
| 7439-92-1 | Lead | 1.5 | U | | P |
| 7439-95-4 | Magnesium | 30000 | | | P |
| 7439-96-5 | Manganese | 69.5 | | E | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 1.7 | B | | P |
| 7440-09-7 | Potassium | 2060 | B | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-23-5 | Sodium | 20700 | | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-62-2 | Vanadium | 1.6 | B | | P |
| 7440-66-6 | Zinc | 0.6 | U | | P |

Color Before: COLORLESS Clarity Before: CLEAR Texture: _____
Color After: COLORLESS Clarity After: CLEAR Artifacts: _____
Comments: _____

U.S. EPA - CLP
1
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKGW07R1009 (DISS)

Lab Name: PROJ AAH GCAL Contract: _____
 Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: _____
 Matrix: (soil / water) Water Lab Sample ID: 20403220809
 Level: (low / med) _____ Date Received: 03/19/04
 % Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|--------------|----|
| 7429-90-5 | Aluminum | 25.8 | U | | P |
| 7440-36-0 | Antimony | 3.7 | U | | P |
| 7440-38-2 | Arsenic | 2.9 | U | | P |
| 7440-39-3 | Barium | 113 | B | | P |
| 7440-41-7 | Beryllium | 0.1 | B | | P |
| 7440-43-9 | Cadmium | 0.2 | U | | P |
| 7440-70-2 | Calcium | 185000 | | | P |
| 7440-47-3 | Chromium | 1.1 | B | | P |
| 7440-48-4 | Cobalt | 0.7 | B | | P |
| 7440-50-8 | Copper | 1.2 | U | | P |
| 7439-89-6 | Iron | 32.9 | B | E | P |
| 7439-92-1 | Lead | 1.5 | U | | P |
| 7439-95-4 | Magnesium | 26300 | | | P |
| 7439-96-5 | Manganese | 914 | | E | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 2.2 | B | | P |
| 7440-09-7 | Potassium | 2350 | B | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-23-5 | Sodium | 25200 | | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-62-2 | Vanadium | 0.8 | U | | P |
| 7440-66-6 | Zinc | 0.6 | U | | P |

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Color Before: COLORLESS Clarity Before: CLEAR Texture: _____
 Color After: COLORLESS Clarity After: CLEAR Artifacts: _____
 Comments: _____

U.S. EPA - CLP
1
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKGW591009 (DISS)

Lab Name: PROJ AAH GCAL Contract: _____
 Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: _____
 Matrix: (soil / water) Water Lab Sample ID: 20403220810
 Level: (low / med) _____ Date Received: 03/19/04
 % Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|--------------|----|
| 7429-90-5 | Aluminum | 27.2 | B | | P |
| 7440-36-0 | Antimony | 3.7 | U | | P |
| 7440-38-2 | Arsenic | 2.9 | U | | P |
| 7440-39-3 | Barium | 21.8 | B | | P |
| 7440-41-7 | Beryllium | 0.1 | B | | P |
| 7440-43-9 | Cadmium | 0.2 | U | | P |
| 7440-70-2 | Calcium | 239000 | | | P |
| 7440-47-3 | Chromium | 1.8 | B | | P |
| 7440-48-4 | Cobalt | 0.4 | B | | P |
| 7440-50-8 | Copper | 2.1 | B | | P |
| 7439-89-6 | Iron | 28.8 | B | E | P |
| 7439-92-1 | Lead | 1.5 | U | | P |
| 7439-95-4 | Magnesium | 49000 | | | P |
| 7439-96-5 | Manganese | 4.5 | B | E | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 2.6 | B | | P |
| 7440-09-7 | Potassium | 32800 | | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-23-5 | Sodium | 166000 | | | P |
| 7440-28-0 | Thallium | 3.1 | B | | P |
| 7440-62-2 | Vanadium | 1.4 | B | | P |
| 7440-66-6 | Zinc | 3.1 | B | | P |

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Color Before: COLORLESS Clarity Before: CLEAR Texture: _____
 Co or After: COLORLESS Clarity After: CLEAR Artifacts: _____
 Comments: _____

U.S. EPA - CLP
1
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKGW62A1009 (DISS)

Lab Name: PROJ AAH GCAL

Contract: _____

Lab Code: LA024

Case No.: _____

SAS No.: _____

SDG No.: _____

Matrix: (soil / water) Water

Lab Sample ID: 20403220812

Level: (low / med) _____

Date Received: 03/19/04

% Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum | 25.8 | U | | P |
| 7440-36-0 | Antimony | 3.7 | U | | P |
| 7440-38-2 | Arsenic | 2.9 | U | | P |
| 7440-39-3 | Barium | 111 | B | | P |
| 7440-41-7 | Beryllium | 0.1 | U | | P |
| 7440-43-9 | Cadmium | 0.2 | U | | P |
| 7440-70-2 | Calcium | 122000 | | | P |
| 7440-47-3 | Chromium | 2.1 | B | | P |
| 7440-48-4 | Cobalt | 0.5 | B | | P |
| 7440-50-8 | Copper | 1.2 | U | | P |
| 7439-89-6 | Iron | 14.1 | U | | P |
| 7439-92-1 | Lead | 1.5 | U | | P |
| 7439-95-4 | Magnesium | 48700 | | | P |
| 7439-96-5 | Manganese | 164 | | E | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 1.8 | B | | P |
| 7440-09-7 | Potassium | 11100 | | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-23-5 | Sodium | 120000 | | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-62-2 | Vanadium | 1.6 | B | | P |
| 7440-66-6 | Zinc | 0.6 | U | | P |

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Color Before: COLORLESS

Clarity Before: CLEAR

Texture: _____

Color After: COLORLESS

Clarity After: CLEAR

Artifacts: _____

Comments:

U.S. EPA - CLP
1
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKGW641009 (DISS)

Lab Name: PROJ AAH GCAL

Contract: _____

Lab Code: LA024

Case No.: _____

SAS No.: _____

SDG No.: _____

Matrix: (soil / water) Water

Lab Sample ID: 20403220813

Level: (low / med) _____

Date Received: 03/19/04

% Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|--------------|----|
| 7429-90-5 | Aluminum | 25.8 | U | | P |
| 7440-36-0 | Antimony | 3.7 | U | | P |
| 7440-38-2 | Arsenic | 2.9 | U | | P |
| 7440-39-3 | Barium | 28.3 | B | | P |
| 7440-41-7 | Beryllium | 0.1 | B | | P |
| 7440-43-9 | Cadmium | 0.2 | U | | P |
| 7440-70-2 | Calcium | 176000 | | | P |
| 7440-47-3 | Chromium | 1.7 | B | | P |
| 7440-48-4 | Cobalt | 1.8 | B | | P |
| 7440-50-8 | Copper | 1.2 | U | | P |
| 7439-89-6 | Iron | 14.1 | U | E | P |
| 7439-92-1 | Lead | 1.5 | U | | P |
| 7439-95-4 | Magnesium | 56700 | | | P |
| 7439-96-5 | Manganese | 1170 | | E | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 7.8 | B | | P |
| 7440-09-7 | Potassium | 12900 | | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-23-5 | Sodium | 53900 | | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-62-2 | Vanadium | 0.8 | U | | P |
| 7440-66-6 | Zinc | 0.6 | U | | P |

Color Before: COLORLESS

Clarity Before: CLEAR

Texture: _____

Color After: COLORLESS

Clarity After: CLEAR

Artifacts: _____

Comments:

INITIAL AND CONTINUING CALIBRATION VERIFICATION

Lab Name: GCAL

Contract: _____

Lab Code: LA024 Case No.: _____

SAS No.: _____ SDG No.: 204031909

Calibration Source: 106-61-2 CPI

Instrument ID: ICP5 ICAL ID: 2

Date Analyzed: 03/25/04 Time: 1050

CRDL STANDARD

| <i>Analyte</i> | <i>True</i> | <i>Found</i> | <i>CAL %R</i> | <i>Units</i> | <i>Method</i> | <i>Type</i> |
|----------------|-------------|--------------|---------------|--------------|----------------------|-------------|
| Antimony | 120 | 123 | 103 | ug/L | ILM04.1 - CLP Metals | P |
| Arsenic | 20.0 | 24.7 | 123.5 | ug/L | ILM04.1 - CLP Metals | P |
| Beryllium | 10.0 | 9.60 | 96 | ug/L | ILM04.1 - CLP Metals | P |
| Cadmium | 10.0 | 9.40 | 94 | ug/L | ILM04.1 - CLP Metals | P |
| Chromium | 20.0 | 19.9 | 99 | ug/L | ILM04.1 - CLP Metals | P |
| Cobalt | 100 | 95.2 | 95 | ug/L | ILM04.1 - CLP Metals | P |
| Copper | 50.0 | 42.0 | 84 | ug/L | ILM04.1 - CLP Metals | P |
| Lead | 6.00 | 5.80 | 97 | ug/L | ILM04.1 - CLP Metals | P |
| Manganese | 30.0 | 29.2 | 97 | ug/L | ILM04.1 - CLP Metals | P |
| Nickel | 80.0 | 78.0 | 97 | ug/L | ILM04.1 - CLP Metals | P |
| Selenium | 10.0 | 12.2 | 122 | ug/L | ILM04.1 - CLP Metals | P |
| Silver | 20.0 | 19.6 | 98 | ug/L | ILM04.1 - CLP Metals | P |
| Thallium | 20.0 | 25.2 | 126 | ug/L | ILM04.1 - CLP Metals | P |
| Vanadium | 100 | 97.2 | 97 | ug/L | ILM04.1 - CLP Metals | P |
| Zinc | 40.0 | 29.5 | 74 | ug/L | ILM04.1 - CLP Metals | P |

INITIAL AND CONTINUING CALIBRATION VERIFICATION

Lab Name: GCAL

Contract: _____

Lab Code: LA024 Case No.: _____

SAS No.: _____ SDG No.: 204031909

Calibration Source: 106-61-2 CPI

Instrument ID: ICP5 ICAL ID: 2

Date Analyzed: 03/25/04 Time: 1340

CRDL STANDARD

| Analyte | True | Found | CAL %R | Units | Method | Type |
|-----------|------|-------|--------|-------|----------------------|------|
| Antimony | 120 | 124 | 103 | ug/L | ILM04.1 - CLP Metals | P |
| Arsenic | 20.0 | 21.2 | 106 | ug/L | ILM04.1 - CLP Metals | P |
| Beryllium | 10.0 | 9.60 | 96 | ug/L | ILM04.1 - CLP Metals | P |
| Cadmium | 10.0 | 9.30 | 93 | ug/L | ILM04.1 - CLP Metals | P |
| Chromium | 20.0 | 20.2 | 101 | ug/L | ILM04.1 - CLP Metals | P |
| Cobalt | 100 | 94.0 | 94 | ug/L | ILM04.1 - CLP Metals | P |
| Copper | 50.0 | 40.7 | 81 | ug/L | ILM04.1 - CLP Metals | P |
| Lead | 6.00 | 5.80 | 96 | ug/L | ILM04.1 - CLP Metals | P |
| Manganese | 30.0 | 29.5 | 98 | ug/L | ILM04.1 - CLP Metals | P |
| Nickel | 80.0 | 77.5 | 97 | ug/L | ILM04.1 - CLP Metals | P |
| Selenium | 10.0 | 10.0 | 100 | ug/L | ILM04.1 - CLP Metals | P |
| Silver | 20.0 | 19.2 | 96 | ug/L | ILM04.1 - CLP Metals | P |
| Thallium | 20.0 | 29.3 | (147) | ug/L | ILM04.1 - CLP Metals | P |
| Vanadium | 100 | 98.4 | 98 | ug/L | ILM04.1 - CLP Metals | P |
| Zinc | 40.0 | 28.6 | (71) | ug/L | ILM04.1 - CLP Metals | P |

INITIAL AND CONTINUING CALIBRATION VERIFICATION

Lab Name: GCAL

Contract: _____

Lab Code: LA024 Case No.: _____

SAS No.: _____ SDG No.: 204031909

Calibration Source: 106-61-2 CPI

Instrument ID: ICP5 ICAL ID: 2

Date Analyzed: 03/25/04 Time: 1635

CRDL STANDARD

| Analyte | True | Found | CAL %R | Units | Method | Type |
|-----------|------|-------|--------|-------|----------------------|------|
| Antimony | 120 | 125 | 104 | ug/L | ILM04.1 - CLP Metals | P |
| Arsenic | 20.0 | 24.0 | 120 | ug/L | ILM04.1 - CLP Metals | P |
| Beryllium | 10.0 | 9.50 | 95 | ug/L | ILM04.1 - CLP Metals | P |
| Cadmium | 10.0 | 9.50 | 95 | ug/L | ILM04.1 - CLP Metals | P |
| Chromium | 20.0 | 19.7 | 98 | ug/L | ILM04.1 - CLP Metals | P |
| Cobalt | 100 | 95.2 | 95 | ug/L | ILM04.1 - CLP Metals | P |
| Copper | 50.0 | 39.5 | 79 | ug/L | ILM04.1 - CLP Metals | P |
| Lead | 6.00 | 6.10 | 102 | ug/L | ILM04.1 - CLP Metals | P |
| Manganese | 30.0 | 29.2 | 97 | ug/L | ILM04.1 - CLP Metals | P |
| Nickel | 80.0 | 77.9 | 97 | ug/L | ILM04.1 - CLP Metals | P |
| Selenium | 10.0 | 13.9 | 139 | ug/L | ILM04.1 - CLP Metals | P |
| Silver | 20.0 | 18.8 | 94 | ug/L | ILM04.1 - CLP Metals | P |
| Thallium | 20.0 | 26.1 | 130 | ug/L | ILM04.1 - CLP Metals | P |
| Vanadium | 100 | 97.2 | 97 | ug/L | ILM04.1 - CLP Metals | P |
| Zinc | 40.0 | 27.6 | 69 | ug/L | ILM04.1 - CLP Metals | P |

U.S. EPA - CLP

3

BLANKS

Lab Name: PROJ AAH GCAL

Contract: _____

Lab Code: LA024

Case No.: _____

SAS No.: _____

SDG No. _____

Preparation Blank Matrix: (soil / water) WaterPreparation Blank Concentration Units: (ug/L / mg/kg) ug/L

| Analyte | Initial Calib. Blank (ug/L) | | Continuing Calibration Blank (ug/L) | | | | | | Preparation Blank | | M |
|-----------|-----------------------------|---|-------------------------------------|---|------|---|------|---|-------------------|---|----|
| | (ug/L) | C | 1 | C | 2 | C | 3 | C | C | | |
| Aluminum | 25.8 | U | 25.8 | U | 25.8 | U | 25.8 | U | 25.800 | U | P |
| Antimony | 9.2 | B | 7.7 | B | 3.7 | U | 3.9 | B | 4.334 | B | P |
| Arsenic | 2.9 | U | 2.9 | U | 4.5 | B | 4.5 | B | 2.900 | U | P |
| Barium | 0.3 | U | 0.3 | U | 0.3 | U | 0.3 | U | 0.604 | B | P |
| Beryllium | 0.1 | U | 0.1 | U | 0.1 | U | 0.1 | U | 0.100 | U | P |
| Cadmium | 1.3 | B | 1.6 | B | 0.2 | U | 2.8 | B | 0.470 | B | P |
| Calcium | 7.5 | U | 7.5 | U | 9.8 | B | 7.5 | U | 9.654 | B | P |
| Chromium | 0.8 | U | 0.8 | U | 0.8 | U | 0.8 | U | 0.800 | U | P |
| Cobalt | 0.4 | U | 0.4 | U | 0.4 | U | 0.6 | B | 0.793 | B | P |
| Copper | 16.7 | B | 19.8 | B | 4.5 | B | 12.8 | B | 20.742 | B | P |
| Iron | 14.1 | U | 14.1 | U | 14.1 | U | 14.1 | U | 17.689 | B | P |
| Lead | 2.2 | B | 1.5 | U | 1.5 | U | 1.5 | U | 1.500 | U | P |
| Magnesium | 36.7 | U | 36.7 | U | 36.7 | U | 81.2 | B | 36.700 | U | P |
| Manganese | 0.2 | U | 0.2 | U | 0.2 | U | 0.2 | U | 0.200 | U | P |
| Mercury | 0.1 | U | 0.1 | U | 0.1 | U | 0.1 | U | 0.100 | U | AV |
| Nickel | 0.7 | U | 0.7 | U | 0.7 | U | 1.2 | B | 1.140 | B | P |
| Potassium | 42.1 | U | 49.4 | B | 42.1 | U | 42.1 | U | 42.100 | U | P |
| Selenium | 4.4 | U | 4.4 | U | 4.4 | U | 4.4 | U | 4.400 | B | P |
| Silver | 0.4 | U | 0.4 | U | -0.6 | B | -0.7 | B | 0.400 | U | P |
| Sodium | 45.4 | U | 45.4 | U | 45.4 | U | 45.4 | U | 89.366 | B | P |
| Thallium | 4.6 | B | 8.0 | B | 2.6 | U | 4.4 | B | 3.425 | B | P |
| Vanadium | 0.8 | U | 0.9 | B | 1.2 | B | 0.9 | B | 1.283 | B | P |
| Zinc | 0.6 | U | 2.5 | B | -4.5 | B | 0.6 | U | 11.747 | B | P |
| Cyanide | 0.5 | U | 0.5 | U | 0.5 | U | 0.5 | U | 0.500 | U | AS |

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3

BLANKS

Lab Name: PROJ AAH GCAL

Contract: _____

Lab Code: LA024 Case No.: _____

SAS No.: _____ SDG No.: _____

Preparation Blank Matrix: (soil / water) WaterPreparation Blank Concentration Units: (ug/L / mg/kg) ug/L

| Analyte | Initial Calib. Blank (ug/L) | Continuing Calibration Blank (ug/L) | | | | | | Prepa- ration Blank C | M |
|-----------|--------------------------------------|-------------------------------------|---|-------|---|---|---|--------------------------------|----|
| | | 4 | C | 5 | C | 6 | C | | |
| Aluminum | | 25.8 | U | 25.8 | U | | | 25.800 | P |
| Antimony | | 3.7 | U | 5.8 | B | | | 3.700 | P |
| Arsenic | | 2.9 | U | 3.3 | B | | | 2.900 | P |
| Barium | | 0.3 | U | 0.3 | U | | | 0.300 | P |
| Beryllium | | 0.1 | U | 0.1 | U | | | 0.100 | P |
| Cadmium | | 0.2 | U | 2.4 | B | | | 0.785 | P |
| Calcium | | 7.5 | U | 7.5 | U | | | 15.883 | P |
| Chromium | | 0.8 | U | 0.8 | U | | | 0.800 | P |
| Cobalt | | 0.4 | U | 0.5 | B | | | 0.859 | P |
| Copper | | 1.4 | B | 11.1 | B | | | 9.684 | P |
| Iron | | 14.1 | U | 14.1 | U | | | 17.257 | P |
| Lead | | 1.5 | U | 1.5 | U | | | 1.500 | P |
| Magnesium | | 36.7 | U | 116.7 | B | | | 57.052 | P |
| Manganese | | 0.2 | U | 0.2 | U | | | 0.200 | P |
| Mercury | | 0.1 | U | | | | | 0.100 | AV |
| Nickel | | 0.7 | U | 0.7 | U | | | 1.690 | P |
| Potassium | | 42.1 | U | 42.1 | U | | | 42.100 | P |
| Selenium | | 4.4 | U | 4.4 | U | | | 4.400 | P |
| Silver | | 0.4 | U | -0.4 | B | | | -0.593 | P |
| Sodium | | 45.4 | U | 45.4 | U | | | 69.029 | P |
| Thallium | | 3.3 | B | 4.5 | B | | | 3.384 | P |
| Vanadium | | 0.8 | U | 0.8 | U | | | 0.800 | P |
| Zinc | | -6.7 | B | -2.2 | B | | | 8.8059 | P |
| Cyanide | | | | | | | | | AS |

MS/MSD RECOVERY

Lab Name: PROJ AAH GCAL

Contract: _____

Lab Code: LA024 Case No.: _____

SAS No.: _____ SDG No.: _____

Matrix Spike - EPA Sample No: SKGW611009SAMPLE NO. : 20403190905

| COMPOUND | UNITS | SPIKE ADDED | SAMPLE CONCENTRATION | MS CONCENTRATION | MS % REC | # | QC. LIMITS |
|-----------|-------|----------------|-------------------------|---------------------|-------------|---|------------|
| Aluminum | ug/L | 2000 | 452 | 2520 | 103 | | 75 - 125 |
| Antimony | ug/L | 100 | 4.8 | 116 | 111 | | 75 - 125 |
| Arsenic | ug/L | 40 | 2.9 | 50 | 125 | | 75 - 125 |
| Barium | ug/L | 2000 | 44.1 | 2120 | 104 | | 75 - 125 |
| Beryllium | ug/L | 50 | .2 | 53.8 | 107 | | 75 - 125 |
| Cadmium | ug/L | 50 | .3 | 50.4 | 100 | | 75 - 125 |
| Chromium | ug/L | 200 | 1.9 | 209 | 104 | | 75 - 125 |
| Cobalt | ug/L | 500 | 1.7 | 498 | 99 | | 75 - 125 |
| Copper | ug/L | 250 | 22.2 | 279 | 103 | | 75 - 125 |
| Iron | ug/L | 1000 | 2430 | 3390 | 95 | | 75 - 125 |
| Lead | ug/L | 20 | 22.1 | 42.2 | 100 | | 75 - 125 |
| Manganese | ug/L | 500 | 527 | 1060 | 106 | | 75 - 125 |
| Mercury | ug/L | 5 | .1 | 5.1 | 103 | | 75 - 125 |
| Nickel | ug/L | 500 | 4.3 | 503 | 100 | | 75 - 125 |
| Selenium | ug/L | 10 | 4.4 | 4.5 | 45 | N | 75 - 125 |
| Silver | ug/L | 50 | .4 | 51.5 | 103 | | 75 - 125 |
| Thallium | ug/L | 50 | 2.6 | 48.3 | 97 | | 75 - 125 |
| Vanadium | ug/L | 500 | 2.1 | 531 | 106 | | 75 - 125 |
| Zinc | ug/L | 500 | 7.3 | 480 | 95 | | 75 - 125 |
| Cyanide | ug/L | 100 | .5 | 86.1 | 86 | | 75 - 125 |

Column to be used to flag recovery and RPD values with an asterisk

* Values outside of QC limits

RPD : 0 out of 0 outside limitsSpike Recovery: 1 out of 20 outside limits

MS/MSD RECOVERY

Lab Name: PROJ AAH GCAL

Contract: _____

Lab Code: LA024

Case No.: _____

SAS No.: _____

SDG No.: _____

Matrix Spike - EPA Sample No: SKGW611009(DISS)

SAMPLE NO. : 20403190914

| COMPOUND | UNITS | SPIKE ADDED | SAMPLE CONCENTRATION | MS CONCENTRATION | MS % REC | # | QC. LIMITS |
|-----------|-------|----------------|-------------------------|---------------------|-------------|---|------------|
| Aluminum | ug/L | 2000 | 25.8 | 2150 | 107 | | 75 - 125 |
| Antimony | ug/L | 100 | 4.5 | 117 | 113 | | 75 - 125 |
| Arsenic | ug/L | 40 | 2.9 | 48.4 | 121 | | 75 - 125 |
| Barium | ug/L | 2000 | 39.4 | 2110 | 104 | | 75 - 125 |
| Beryllium | ug/L | 50 | .1 | 54.3 | 108 | | 75 - 125 |
| Cadmium | ug/L | 50 | .3 | 50.1 | 100 | | 75 - 125 |
| Chromium | ug/L | 200 | 1.1 | 208 | 104 | | 75 - 125 |
| Cobalt | ug/L | 500 | 1.4 | 489 | 98 | | 75 - 125 |
| Copper | ug/L | 250 | 8 | 272 | 106 | | 75 - 125 |
| Iron | ug/L | 1000 | 187 | 1210 | 103 | | 75 - 125 |
| Lead | ug/L | 20 | 1.5 | 18.7 | 93 | | 75 - 125 |
| Manganese | ug/L | 500 | 485 | 991 | 101 | | 75 - 125 |
| Mercury | ug/L | 5 | .1 | 4.1 | 81 | | 75 - 125 |
| Nickel | ug/L | 500 | 4.2 | 491 | 97 | | 75 - 125 |
| Selenium | ug/L | 10 | 4.4 | 7.3 | 73 | N | 75 - 125 |
| Silver | ug/L | 50 | .4 | 52.9 | 106 | | 75 - 125 |
| Thallium | ug/L | 50 | 2.6 | 53.5 | 107 | | 75 - 125 |
| Vanadium | ug/L | 500 | 1.2 | 532 | 106 | | 75 - 125 |
| Zinc | ug/L | 500 | .6 | 467 | 93 | | 75 - 125 |

Column to be used to flag recovery and RPD values with an asterisk

* Values outside of QC limits

RPD : 0 out of 0 outside limits

Spike Recovery: 1 out of 19 outside limits

FORM V (PART 1) - IN

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U.S. EPA - CLP
5B
POST DIGEST SPIKE SAMPLE RECOVERY

EPA SAMPLE NO.

SKGW611009PDS

Lab Name: PROJ AAH GCAL

Lab Code: LA024

Case No.: _____

Contract: _____

Matrix: (soil / water) Water

SAS No.: _____

SDG No.: _____

% Solids for Sample: _____

Level: (low / med) _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| Analyte | Control Limit | | Spiked Sample | | Sample | | Spike Added (SA) | % R | Q | M |
|-----------|---------------|--|---------------|---|-------------|---|------------------|-----|---|---|
| | %R | | Result (SSR) | C | Result (SR) | C | | | | |
| Aluminum | | | 527 | | 25.8 | U | | 0 | | P |
| Antimony | | | 142 | | 4.8 | B | 120 | 114 | | P |
| Arsenic | | | 27.2 | | 2.9 | U | 20 | 136 | | P |
| Barium | | | 44.2 | B | .3 | U | | 0 | | P |
| Beryllium | | | 11 | | .2 | B | 10 | 109 | | P |
| Cadmium | | | 10.2 | | .3 | B | 10 | 99 | | P |
| Calcium | | | 185000 | | 7.5 | U | | 0 | | P |
| Chromium | | | 22.7 | | 1.9 | B | 20 | 104 | | P |
| Cobalt | | | 101 | | 1.7 | B | 100 | 100 | | P |
| Copper | | | 69.1 | | 22.2 | B | 50 | 94 | | P |
| Iron | | | 2400 | | 14.1 | U | | 0 | | P |
| Lead | | | 28 | | 22.1 | | 6 | 97 | | P |
| Magnesium | | | 29300 | | 36.7 | U | | 0 | | P |
| Manganese | | | 554 | | 527 | | 30 | 90 | | P |
| Nickel | | | 82.6 | | 4.3 | B | 80 | 98 | | P |
| Potassium | | | 6960 | | 42.1 | U | | 0 | | P |
| Selenium | | | 4.9 | B | 4.4 | U | 10 | 49 | | P |
| Silver | | | 20.3 | | .4 | U | 20 | 101 | | P |
| Sodium | | | 27000 | | 45.4 | U | | 0 | | P |
| Thallium | | | 17.4 | | 2.6 | U | 20 | 87 | | P |
| Vanadium | | | 110 | | 2.1 | B | 100 | 108 | | P |
| Zinc | | | 44.3 | | 7.3 | B | 40 | 93 | | P |

Comments:

U.S. EPA - CLP
5B
POST DIGEST SPIKE SAMPLE RECOVERY

EPA SAMPLE NO.

SKGW611009(DISS)PDS

Lab Name: PROJ AAH GCAL

Lab Code: LA024

Case No.: _____

Contract: _____

Matrix: (soil / water) Water

SAS No.: _____

SDG No.: _____

% Solids for Sample: _____

Level: (low / med) _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| Analyte | Control Limit %R | Spiked Sample Result (SSR) | C | Sample Result (SR) | C | Spike Added (SA) | % R | Q | M |
|-----------|------------------------|----------------------------------|---|-----------------------|---|---------------------|-----|---|---|
| Aluminum | | 25.8 | U | 25.8 | U | | 0 | | P |
| Antimony | | 136 | | 4.5 | B | 120 | 110 | | P |
| Arsenic | | 26.1 | | 2.9 | U | 20 | 131 | | P |
| Barium | | 39.9 | B | .3 | U | | 0 | | P |
| Beryllium | | 10.9 | | .1 | B | 10 | 107 | | P |
| Cadmium | | 10 | | .3 | B | 10 | 97 | | P |
| Calcium | | 191000 | | 7.5 | U | | 0 | | P |
| Chromium | | 21.6 | | 1.1 | B | 20 | 103 | | P |
| Cobalt | | 99 | | 1.4 | B | 100 | 98 | | P |
| Copper | | 60.1 | | 8 | B | 50 | 104 | | P |
| Iron | | 181 | | 14.1 | U | | 0 | | P |
| Lead | | 4 | | 1.5 | U | 6 | 67 | | P |
| Magnesium | | 29000 | | 36.7 | U | | 0 | | P |
| Manganese | | 524 | | 485 | | 30 | 128 | | P |
| Nickel | | 78.2 | | 4.2 | B | 80 | 93 | | P |
| Potassium | | 6970 | | 42.1 | U | | 0 | | P |
| Selenium | | 4.4 | U | 4.4 | U | 10 | 0 | | P |
| Silver | | 20.3 | | .4 | U | 20 | 102 | | P |
| Sodium | | 28000 | | 45.4 | U | | 0 | | P |
| Thallium | | 19.8 | | 2.6 | U | 20 | 99 | | P |
| Vanadium | | 108 | | 1.2 | B | 100 | 107 | | P |
| Zinc | | 40.8 | | .6 | U | 40 | 102 | | P |

Comments:

U.S. EPA - CLP

6

DUPLICATES

EPA SAMPLE NO

SKGW61DUP1009

Lab Name: PROJ AAH GCAL

Lab Code: LA024

Case No.:

Contract:

Matrix: (soil / water) Water

SAS No.:

SDG No.:

Level: (low / med)

% Solids for Sample:

% Solids for Duplicate:

Concentration Units (ug/L or mg/kg dry weight) ug/L

| Analyte | Control Limit | Sample (S) | C | Duplicate (D) | C | RPD | Q | M |
|-----------|---------------|------------|---|---------------|---|-----|---|----|
| Alum num | 0 - 200 | 452 | | 470 | | 18 | | P |
| Antimony | - | 4.8 | B | 5.5 | B | 14 | | P |
| Arsenic | - | 2.9 | U | 3.7 | B | 200 | | P |
| Barium | - | 44.1 | B | 45.4 | B | 3 | | P |
| Beryllium | - | .2 | B | .2 | B | 0 | | P |
| Cadmium | - | .3 | B | .3 | B | 0 | | P |
| Calcium | 0 - 20 | 187000 | | 192000 | | 3 | | P |
| Chromium | - | 1.9 | B | 1.6 | B | 17 | | P |
| Cobalt | - | 1.7 | B | 1.6 | B | 6 | | P |
| Copper | - | 22.2 | B | 18.6 | B | 18 | | P |
| Iron | 0 - 20 | 2430 | | 2490 | | 2 | | P |
| Lead | 0 - 20 | 22.1 | | 22.5 | | 2 | | P |
| Magnesium | 0 - 20 | 30000 | | 30300 | | 1 | | P |
| Manganese | 0 - 20 | 527 | | 530 | | .6 | | P |
| Mercury | - | .1 | U | .1 | U | 0 | | AV |
| Nicke | - | 4.3 | B | 3.7 | B | 15 | | P |
| Potassium | 0 - 5000 | 6950 | | 7140 | | 190 | | P |
| Selen um | - | 4.4 | U | 4.4 | U | 0 | | P |
| Silver | - | .4 | U | .4 | U | 0 | | P |
| Sodium | 0 - 20 | 27000 | | 27800 | | 3 | | P |
| Thallium | - | 2.6 | U | 2.6 | U | 0 | | P |
| Vanadium | - | 2.1 | B | 1.7 | B | 21 | | P |
| Zinc | - | 7.3 | B | 7 | B | 4 | | P |
| Cyanide | - | .5 | U | .8 | B | 91 | | AS |

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RESUBMITTED

U.S. EPA - CLP

6

DUPLICATES

EPA SAMPLE NO

SKGW61DUP1009 (DISS)

Lab Name: PROJ AAH GCALLab Code: LA024

Case No.: _____

Contract: _____

Matrix: (soil / water) Water

SAS No.: _____

SDG No.: _____

Level: (low / med) _____

% Solids for Sample: _____

% Solids for Duplicate: _____

Concentration Units (ug/L or mg/kg dry weight) ug/L

| Analyte | Control Limit | Sample (S) | C | Duplicate (D) | C | RPD | Q | M |
|-----------|---------------|------------|---|---------------|---|-----|---|----|
| Aluminum | - | 25.8 | U | 25.8 | U | 0 | | P |
| Antimony | - | 4.5 | B | 3.7 | U | 200 | | P |
| Arsenic | - | 2.9 | U | 2.9 | U | 0 | | P |
| Barium | - | 39.4 | B | 38.9 | B | 1 | | P |
| Beryllium | - | .1 | B | .1 | B | 0 | | P |
| Cadmium | - | .3 | B | .2 | U | 200 | | P |
| Calcium | 0 - 20 | 191000 | | 186000 | | 3 | | P |
| Chromium | - | 1.1 | B | 1.3 | B | 17 | | P |
| Cobalt | - | 1.4 | B | 1.3 | B | 7 | | P |
| Copper | - | 8 | B | 5.2 | B | 42 | | P |
| Iron | 0 - 100 | 187 | | 195 | | 8 | | P |
| Lead | - | 1.5 | U | 1.5 | U | 0 | | P |
| Magnesium | 0 - 20 | 29100 | | 28400 | | 2 | | P |
| Manganese | 0 - 20 | 485 | | 477 | | 2 | | P |
| Mercury | - | .1 | U | .1 | U | 0 | | AV |
| Nickel | - | 4.2 | B | 3.6 | B | 15 | | P |
| Potassium | 0 - 5000 | 6990 | | 6860 | | 130 | | P |
| Selenium | - | 4.4 | U | 4.4 | U | 0 | | P |
| Silver | - | .4 | U | .4 | U | 0 | | P |
| Sodium | 0 - 20 | 27900 | | 27200 | | 3 | | P |
| Thallium | - | 2.6 | U | 5.6 | B | 200 | | P |
| Vanadium | - | 1.2 | B | 1.8 | B | 40 | | P |
| Zinc | - | .6 | U | .6 | U | 0 | | P |

U.S. EPA - CLP
7
LABORATORY CONTROL SAMPLE

Lab Name: PROJ AAH GCAL Contract: _____
 Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: _____
 Solid LCS Source: _____
 Aqueous LCS Source: 310006 HIGH PURITY~323003 HIGH

| Analyte | Aqueous (ug/L) | | | Solid (mg/kg) | | | |
|-----------|----------------|-------|-----|---------------|-------|---|-----|
| | True | Found | % R | True | Found | C | % R |
| Aluminum | 2000 | 2190 | 109 | | | | |
| Antimony | 500 | 557 | 111 | | | | |
| Arsenic | 2000 | 2090 | 104 | | | | |
| Barium | 2000 | 2070 | 104 | | | | |
| Beryllium | 50.0 | 54.2 | 108 | | | | |
| Cadmium | 50.0 | 51.5 | 103 | | | | |
| Calcium | 12500 | 13200 | 106 | | | | |
| Chromium | 200 | 206 | 103 | | | | |
| Cobalt | 500 | 501 | 100 | | | | |
| Copper | 250 | 280 | 112 | | | | |
| Iron | 1000 | 1080 | 108 | | | | |
| Lead | 500 | 512 | 102 | | | | |
| Magnesium | 12500 | 12700 | 102 | | | | |
| Manganese | 500 | 521 | 104 | | | | |
| Nickel | 500 | 509 | 102 | | | | |
| Potassium | 12500 | 12900 | 103 | | | | |
| Selenium | 2000 | 2120 | 106 | | | | |
| Silver | 50.0 | 52.6 | 105 | | | | |
| Sodium | 12500 | 13100 | 105 | | | | |
| Thallium | 2000 | 2050 | 102 | | | | |
| Vanadium | 500 | 530 | 106 | | | | |
| Zinc | 500 | 481 | 96 | | | | |

U.S. EPA - CLP
7
LABORATORY CONTROL SAMPLE

Lab Name: PROJ AAH GCAL Contract: _____
 Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: _____
 Solid LCS Source: _____
 Aqueous LCS Source: 310006 HIGH PURITY~323003 HIGH

| Analyte | Aqueous (ug/L) | | | Solid (mg/kg) | | | |
|-----------|----------------|-------|-----|---------------|-------|---|-----|
| | True | Found | % R | True | Found | C | % R |
| Aluminum | 2000 | 2110 | 105 | | | | |
| Antimony | 500 | 552 | 110 | | | | |
| Arsenic | 2000 | 2080 | 104 | | | | |
| Barium | 2000 | 2060 | 103 | | | | |
| Beryllium | 50.0 | 53.8 | 108 | | | | |
| Cadmium | 50.0 | 51.5 | 103 | | | | |
| Calcium | 12500 | 13000 | 104 | | | | |
| Chromium | 200 | 204 | 102 | | | | |
| Cobalt | 500 | 499 | 100 | | | | |
| Copper | 250 | 275 | 110 | | | | |
| Iron | 1000 | 1040 | 104 | | | | |
| Lead | 500 | 512 | 102 | | | | |
| Magnesium | 12500 | 12500 | 100 | | | | |
| Manganese | 500 | 517 | 103 | | | | |
| Nickel | 500 | 507 | 101 | | | | |
| Potassium | 12500 | 12600 | 101 | | | | |
| Selenium | 2000 | 2120 | 106 | | | | |
| Silver | 50.0 | 51.4 | 103 | | | | |
| Sodium | 12500 | 12800 | 103 | | | | |
| Thallium | 2000 | 2040 | 102 | | | | |
| Vanadium | 500 | 525 | 105 | | | | |
| Zinc | 500 | 475 | 95 | | | | |

U.S. EPA - CLP

9

EPA SAMPLE NO.

ICP SERIAL DILUTIONS

SKGW611009SD

Lab Name: PROJ AAH GCALLab Code: LA024

Case No. _____

Contract: _____

Matrix: (soil / water) Water

SAS No.: _____

SDG No.: _____

Level: (low / med) _____

Concentration Units: ug/L

| Analyte | Initial Sample Result (I) | C | Serial Dilution Result (S) | C | % Difference | Q | M |
|-----------|---------------------------------|---|----------------------------------|---|--------------|--------------|---|
| Aluminum | 452 | | 536 | B | 18.6 | E | P |
| Antimony | 4.8 | B | 18.5 | U | 285 | | P |
| Arsenic | 2.9 | U | 17.7 | B | | | P |
| Barium | 44.1 | B | 43.0 | B | 2.5 | | P |
| Beryllium | 0.2 | B | 0.5 | U | 150 | | P |
| Cadmium | 0.3 | B | 1.0 | U | 233 | | P |
| Calcium | 187000 | | 189000 | | 1.1 | | P |
| Chromium | 1.9 | B | 4.4 | B | 132 | | P |
| Cobalt | 1.7 | B | 2.0 | U | 17.6 | | P |
| Copper | 22.2 | B | 6.0 | U | 73 | E | P |
| Iron | 2430 | | 2400 | | 1.2 | E | P |
| Lead | 22.1 | | 13.8 | B | 37.6 | E | P |
| Magnesium | 30000 | | 29300 | | 2.3 | | P |
| Manganese | 527 | | 525 | | .4 | | P |
| Nickel | 4.3 | B | 4.1 | B | 4.7 | | P |
| Potassium | 6950 | | 6380 | B | 8.2 | | P |
| Selenium | 44.0 | U | 22.0 | U | | | P |
| Silver | 4.0 | U | 2.0 | U | | | P |
| Sodium | 27000 | | 25900 | | 4.1 | | P |
| Thallium | 26.0 | U | 13.0 | U | | | P |
| Vanadium | 2.1 | B | 7.2 | B | 243 | | P |
| Zinc | 7.3 | B | 3.0 | U | 58.9 | E | P |

U.S. EPA - CLP
9
ICP SERIAL DILUTIONS

EPA SAMPLE NO.

SKGW611009(DISS)SD

Lab Name: PROJ AAH GCAL

Lab Code: LA024

Case No.

Contract:

Matrix: (soil / water) Water

SAS No.:

SDG No.:

Level: (low / med)

Concentration Units: ug/L

| Analyte | Initial Sample Result (I) | C | Serial Dilution Result (S) | C | % Difference | Q | M |
|-----------|---------------------------------|---|----------------------------------|---|--------------|---|---|
| Aluminum | 258 | U | 129 | U | | | P |
| Antimony | 4.5 | B | 18.5 | U | 311 | | P |
| Arsenic | 29.0 | U | 14.5 | U | | | P |
| Barium | 39.4 | B | 36.9 | B | 6.3 | | P |
| Beryllium | 0.1 | B | 0.5 | U | 400 | | P |
| Cadmium | 0.3 | B | 1.0 | U | 233 | | P |
| Calcium | 191000 | | 194000 | | 1.6 | | P |
| Chromium | 1.1 | B | 4.0 | U | 264 | | P |
| Cobalt | 1.4 | B | 2.0 | U | 42.9 | | P |
| Copper | 8.0 | B | 6.0 | U | 25 | | P |
| Iron | 187 | | 149 | B | 20.3 | E | P |
| Lead | 15.0 | U | 7.5 | U | | | P |
| Magnesium | 29100 | | 28900 | | .7 | | P |
| Manganese | 485 | | 545 | | 12.4 | E | P |
| Nickel | 4.2 | B | 5.3 | B | 26.2 | | P |
| Potassium | 6990 | | 6620 | B | 5.3 | | P |
| Selenium | 44.0 | U | 22.0 | U | | | P |
| Silver | 4.0 | U | 2.0 | U | | | P |
| Sodium | 27900 | | 27000 | | 3.2 | | P |
| Thallium | 26.0 | U | 13.0 | U | | | P |
| Vanadium | 1.2 | B | 4.0 | U | 233 | | P |
| Zinc | 6.0 | U | 3.0 | U | | | P |

U.S. EPA - CLP

10

INSTRUMENT DETECTION LIMITS (QUARTERLY)

Lab Name: PROJ AAH GCALLab Code: LA024

Case No.: _____

SAS No.: _____

SDG No.: _____

Instrument ID: ICP5Study Date: 04/30/03

| <i>Analyte</i> | <i>Wavelength (nm)</i> | <i>Background</i> | <i>CRDL (ug/L)</i> | <i>IDL (ug/L)</i> | <i>M</i> |
|----------------|----------------------------|-------------------|------------------------|-----------------------|----------|
| Aluminum | 308.210 | | 200 | 25.8 | P |
| Antimony | 206.830 | | 60 | 3.7 | P |
| Arsenic | 193.700 | | 10 | 2.9 | P |
| Barium | 233.520 | | 200 | .3 | P |
| Beryllium | 313.100 | | 5 | .1 | P |
| Cadmium | 214.430 | | 5 | .2 | P |
| Calcium | 315.880 | | 5000 | 7.5 | P |
| Chromium | 267.710 | | 10 | .8 | P |
| Cobalt | 228.610 | | 50 | .4 | P |
| Copper | 324.750 | | 25 | 1.2 | P |
| Iron | 259.940 | | 100 | 14.1 | P |
| Lead | 220.350 | | 3 | 1.5 | P |
| Magnesium | 279.080 | | 5000 | 36.7 | P |
| Manganese | 257.610 | | 15 | .2 | P |
| Nickel | 231.600 | | 40 | .7 | P |
| Potassium | 766.480 | | 5000 | 42.1 | P |
| Selenium | 196.030 | | 5 | 4.4 | P |
| Silver | 328.060 | | 10 | .4 | P |
| Sodium | 589.580 | | 5000 | 45.4 | P |
| Thallium | 190.800 | | 10 | 2.6 | P |
| Vanadium | 290.880 | | 50 | .8 | P |
| Zinc | 213.860 | | 20 | .6 | P |

U.S. EPA - CLP
14
ANALYSIS RUN LOG

Lab Name: PROJ AAH GCAL
Lab Code: LA024 Case No.: _____
Instrument ID Number: ICP5
Start Date: 03/25/04

Contract: _____
SAS No.: _____ SDG No.: _____
Method Type: P
End Date: 03/25/04

Analyte Symbols

| EPA Sample No. | D/F | Time | % R | Al | Sb | As | Ba | Be | Cd | Ca | Cr | Co | Cu | Fe | Pb | Mg | Mn | Hg | Ni | K | Se | Ag | Na | Tl | V | Zn | Cn |
|----------------|-----|------|-----|----|----|----|----|----|----|----|----|----|----|----|----|----|----|----|----|---|----|----|----|----|---|----|----|
| ICV | 1 | 1030 | | | X | X | | X | X | X | X | X | X | X | X | X | X | | X | | X | X | | X | X | X | |
| ICV2 | 1 | 1036 | | X | | | X | | | | | | | | | | | | | X | | | X | | | | |
| ICB | 1 | 1043 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| CRDL | 1 | 1050 | | | X | X | | X | X | | X | X | X | | X | | X | | X | | X | X | | X | X | X | |
| ICSA | 1 | 1057 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| ICSAB | 1 | 1103 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| CCV | 1 | 1109 | | X | | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| CCV2 | 1 | 1115 | | | X | | | | | | | | | | | | | | | | | | | | | | |
| CCB | 1 | 1122 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| MB154054 | 1 | 1129 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKGW611009 | 1 | 1136 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKGW61DUP1009 | 1 | 1143 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKGW611009SD | 5 | 1149 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKGW581009 | 1 | 1156 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKGW58DUP1009 | 1 | 1203 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKGW631009 | 1 | 1209 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKGW06R1009 | 1 | 1215 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKGW07R1009 | 1 | 1221 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKGW591009 | 1 | 1227 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| CCV | 1 | 1236 | | X | | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| CCV2 | 1 | 1243 | | | X | | | | | | | | | | | | | | | | | | | | | | |
| CCB | 1 | 1250 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKGW601009 | 1 | 1257 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKGW62A1009 | 1 | 1303 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKGW641009 | 1 | 1309 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKGWFB1009 | 1 | 1315 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKGW61MS1009 | 1 | 1322 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |

U.S. EPA - CLP
14
ANALYSIS RUN LOG

Lab Name: PROJ AAH GCAL
Lab Code: LA024 Case No.:
Instrument ID Number: ICP5
Start Date: 03/25/04

Contract:
SAS No.: SDG No.:
Method Type: P
End Date: 03/25/04

| Analyte Symbols | | | | | | | | | | | | | | | | | | | | | | | | | | | |
|----------------------|-----|------|-----|----|----|----|----|----|----|----|----|----|----|----|----|----|----|----|----|---|----|----|----|----|---|----|----|
| EPA Sample No. | D/F | Time | % R | Al | Sb | As | Ba | Be | Cd | Ca | Cr | Co | Cu | Fe | Pb | Mg | Mn | Hg | Ni | K | Se | Ag | Na | Tl | V | Zn | Cn |
| SKGW611009PDS | 1 | 1328 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| LCS154055 | 1 | 1334 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| CRDL | 1 | 1340 | | | X | X | | X | X | | X | X | X | | X | | X | | X | | X | X | | X | X | X | |
| ICSA | 1 | 1347 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| ICSAB | 1 | 1353 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| CCV | 1 | 1359 | | X | | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| CCV2 | 1 | 1405 | | | X | | | | | | | | | | | | | | | | | | | | | | |
| CCB | 1 | 1412 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| MB154057 | 1 | 1420 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKGW611009(DISS) | 1 | 1427 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKGW61DUP1009 (DISS) | 1 | 1434 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKGW611009(DISS)SD | 5 | 1441 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKGW581009(DISS) | 1 | 1448 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKGW58DUP1009(DISS) | 1 | 1455 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKGW631009(DISS) | 1 | 1502 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKGW06R1009 (DISS) | 1 | 1509 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKGW07R1009 (DISS) | 1 | 1516 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKGW591009 (DISS) | 1 | 1522 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| CCV | 1 | 1529 | | X | | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| CCV2 | 1 | 1535 | | | X | | | | | | | | | | | | | | | | | | | | | | |
| CCB | 1 | 1542 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKGW601009 (DISS) | 1 | 1549 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKGW62A1009 (DISS) | 1 | 1556 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKGW641009 (DISS) | 1 | 1602 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKGWFB1009 (DISS) | 1 | 1609 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKGW61MS1009(DISS) | 1 | 1616 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKGW611009(DISS)PDS | 1 | 1622 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |

000551
RESUBMITTED

U.S. EPA - CLP
14
ANALYSIS RUN LOG

Lab Name: PROJ AAH GCAL Contract: _____
 Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: _____
 Instrument ID Number: ICP5 Method Type: P
 Start Date: 03/25/04 End Date: 03/25/04

Analyte Symbols

| EPA Sample No. | D/F | Time | % R | Al | Sb | As | Ba | Be | Cd | Ca | Cr | Co | Cu | Fe | Pb | Mg | Mn | Hg | Ni | K | Se | Ag | Na | Tl | V | Zn | Cn |
|----------------|-----|------|-----|----|----|----|----|----|----|----|----|----|----|----|----|----|----|----|----|---|----|----|----|----|---|----|----|
| LCS154058 | 1 | 1629 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| CRDL | 1 | 1635 | | | X | X | | X | X | | X | X | X | | X | | X | | X | | X | X | | X | X | X | |
| ICSA | 1 | 1642 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| ICSAB | 1 | 1648 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| CCV | 1 | 1655 | | X | | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| CCV2 | 1 | 1701 | | | X | | | | | | | | | | | | | | | | | | | | | | |
| CCB | 1 | 1708 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |

000552
RESUBMITTED

DATA VALIDATION REPORT
FOR
SKINNER LANDFILL SITE
EARTH TECH: PROJECT NUMBER 54280
LABORATORY REPORT NUMBER 204032408
PROJECT MANAGER: Ron Rolker
Date: May 25, 2004
Revised Date: August 26, 2005
Data Validator: Mark Kromis

APPENDIX C LIST OF ACRONYMS

| | |
|--------|--|
| BFB | Bromofluorobenzene |
| CC | Continuing Calibration |
| CCV | Continuing Calibration Verification |
| CCB | Continuing Calibration Blanks |
| CLP | Contract Laboratory Program |
| CRDL | Contract Required Detection Limit |
| DFTPP | Decafluorotriphenylphosphine |
| GC/MS | Gas Chromatograph/Mass Spectrometer |
| IC | Initial Calibration |
| ICB | Initial Calibration Blank |
| IDL | Instrument Detection Limit |
| ICP | Inductively Coupled Plasma |
| ICS | Interference Check Sample |
| ICV | Initial Calibration Verification |
| ILM | Inorganic Analysis Multi-Media Multi-Concentration |
| INDAM | Individual A Mixture |
| INDBM | Individual B Mixture |
| mg/L | milligrams per liter |
| MS/MSD | Matrix Spike/Matrix Spike Duplicate |
| OLC | Organic Analysis Low Concentration |
| OLM | Organic Analysis Multi-Media Multi-Concentration |
| %D | Percent Difference |
| % RSD | Percent Relative Standard Deviation |
| PB | Preparation Blanks |
| QC | Quality Control |
| RF | Response Factor |
| RPD | Relative Percent Difference |
| RRF | Relative Response Factor |
| SDG | Sample Delivery Group |
| SOW | Statement of Work |
| µg/L | micrograms per liter |
| US EPA | United States Environmental Protection Agency |
| VOC | Volatile Organic Compounds |
| VTSR | Validated Time of Sample Receipt |

CASE NARRATIVE

The Inorganic portion of this report was revised do to the submission of additional data for the metals that were not reported by GCAL in the original data package. The following metals were included in the resubmission: Aluminum, Calcium, Cobalt, Magnesium, Manganese, Potassium, Sodium, and Vanadium.

DATA VALIDATION SUMMARY – SAMPLE DELIVERY GROUP 204032408 INORGANICS

Validation of the inorganics data, as prepared by Gulf Coast Analytical Laboratories (GCAL) for the samples collected from the Skinner Landfill site in March 2004, was conducted by Earth Tech using the National Functional Guidelines for Inorganic Data Review, (US EPA, February, 1994), as appropriate. The results were reported by GCAL under Sample Delivery Group (SDG) 204032408.

| GCAL # | Sample Description |
|-------------|----------------------|
| 20403240801 | SKSWEB1009 |
| 20403240802 | SKSW511009 |
| 20403240803 | SKSW51MS1009 |
| 20403240805 | SKSW51DUP1009 |
| 20403240806 | SKSW521009 |
| 20403240807 | SKSW531009 |
| 20403240808 | SKSWEB1009 (DISS) |
| 20403240809 | SKSW511009 (DISS) |
| 20403240810 | SKSW51MS1009 (DISS) |
| 20403240811 | SKSW51DUP1009 (DISS) |
| 20403240812 | SKSW521009 (DISS) |
| 20403240813 | SKSW531009 (DISS) |

INTRODUCTION

Analyses of metals were performed according to Contract Laboratory Program (CLP)-Inorganic Analysis Multi-media Multi-concentration ILM04.1 Statement of Work (SOW). Results of the sample analyses are reported by the laboratory as either qualified or unqualified. Unqualified results mean that the reported values maybe used without reservation. Various qualifier codes are used by the laboratory to denote specific information regarding the analytical results. The data validation process is intended to evaluate the data on a technical basis. The data package also was subjected to an internal laboratory quality review prior to submission to Earth Tech for data validation.

During the validation process, laboratory-qualified and unqualified data are verified against all available supporting documentation. Based on this evaluation, qualifier codes may be added, deleted or modified by the data user. Final results are therefore, either qualified or unqualified. Validator-qualified results are annotated with the following codes in accordance with the Functional Guidelines:

- U The constituent was analyzed for, but was not detected above the level of the associated analytical reporting limit. The associated value is either the sample quantitation limit or the sample detection limit.
- J The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample.

- UJ The analyte was not detected above the reported sample quantitation limit. However, the reported quantitation limit is approximate and may or may not represent the actual limit of quantitation necessary to accurately and precisely measure the analyte in the sample.
- R The sample results are rejected due to serious deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte cannot be verified.

Details of the inorganics data validation findings and conclusions are provided in the following sections of this report:

1. Holding Times
2. Calibration
 - A. Initial Calibration (IC)
 - B. Continuing Calibration (CC)
3. Blanks
4. Inductively Coupled Plasma (ICP) Interference Check Sample
5. Laboratory Control Sample (LCS)
6. Duplicate Analysis
7. Spike Sample Analysis
8. ICP Serial Dilution
9. System Performance
10. Documentation
11. Overall Assessment

1. HOLDING TIMES

All samples for inorganics analyses were analyzed within the 180-day holding time for preserved aqueous samples. Mercury analyses were conducted within the 28-day holding time for aqueous samples undergoing CLP protocol. Cyanide analyses were conducted within the 14-day holding time. The cooler temperature upon receipt at the laboratory was within the recommended temperature of 4°C +/- 2°C.

2. CALIBRATION

A. Initial Calibration

The percent recoveries for the Initial Calibration Verification (ICV) standard were within Quality Control (QC) limits for all constituents.

B. Continuing Calibration

The percent recoveries for the Continuing Calibration Verification (CCV) standard were within QC limits for all constituents.

3. BLANKS

The Initial Calibration Blank (ICB), Continuing Calibration Blanks (CCB) and Preparation Blanks (PB) were analyzed at the appropriate frequencies. No constituents were detected in the ICB, CCB, and PB blanks above the corresponding Contract Required Detection Limit (CRDL).

4. ICP INTERFERENCE CHECK SAMPLE

Results for the ICP analysis of the Interference Check Sample (ICS) solution AB were within 20% of the true value.

5. LABORATORY CONTROL SAMPLES

Recoveries were within the control limit (80-120%) for all constituents.

6. DUPLICATE ANALYSIS

The laboratory used sample SKSW511009 for the duplicate sample. The Relative Percent Difference (RPD) between the sample and duplicate results for the dissolved fraction were within the acceptance criteria (<20%) for all target compounds. The Relative Percent Difference (RPD) between the sample and duplicate results for the total fraction were within the acceptance criteria (<20%) for all target compounds with the exception of Aluminum, Arsenic, Barium, Calcium, Chromium, Copper, Iron, Lead, Magnesium, Manganese, Nickel, and Zinc. As per the National Functional Guidelines, if the result from a duplicate analysis for a particular analyte falls outside the appropriate fixed control windows; qualify the results for that analyte in all associated samples of the same matrix as estimated (J).

7. SPIKE SAMPLE ANALYSIS

The laboratory used sample SKSW511009 for the matrix spike sample. The MS percent recoveries were within the acceptance criteria (75%-125%) with the exception of Antimony (35%), Arsenic (151%), Selenium (0%), and Zinc (128%) in the total fraction and Arsenic (128%) and Selenium (0%) in the dissolved fraction.

As per the National Functional Guidelines: if the percent recovery is greater than 30% and less than 74% qualify detected results for that analyte with "J" and non-detected results with "UJ". If the percent recovery is less than 30% qualify detected results for that analyte with "J" and non-detected results with "R".

8. ICP SERIAL DILUTION

As noted in the National Functional Guidelines: If the analyte concentration is at least 50 times above the IDL, its serial dilution analysis must then agree within 10% of the original determination after corrected for dilution. The serial dilution is performed to determine whether any significant chemical or physical interference's exist due to matrix effects. The serial dilution percent differences were within the acceptance criteria for all target analytes with the exception of Aluminum, Barium, Chromium, Iron, Magnesium, Nickel, Potassium, and Vanadium in the total fraction. The serial dilution percent differences were within the acceptance criteria for all target analytes with the exception of Manganese in the dissolved fraction. As per the National Functional Guidelines: if the required 10% Difference criteria are not met qualify the associated data as estimated "J".

9. SYSTEM PERFORMANCE

The analytical system appears to have been working well at the time of these analyses, based on the evaluation of the raw data.

10. DOCUMENTATION

GCAL qualified the total metal results for Beryllium, Cadmium, Cobalt, Mercury and Potassium with (*) on the Form 1's to indicate that the RPD between the original results and its duplicate result exceeded the control limit. The RPD was actually within the control limit therefore the data validator crossed out the (*) with a single line and dated and initialed the bottom of the report.

11. OVERALL ASSESSMENT

The percent recoveries for Copper in the Contract Required Detection Limit (CRDL) standards were 80%, 77% and 78%.

If the CRDL is below 80% then detected results are qualified as estimated with "J" and the non-detected results were qualified with "UJ".

The results are acceptable with the validator-added qualifiers.

**DATA VALIDATION SUMMARY – SAMPLE DELIVERY GROUP 204032408
SEMIVOLATILE ORGANICS**

Validation of the Gas Chromatograph/Mass Spectrometer (GC/MS) semi-volatile organics data, as prepared by Gulf Coast Analytical Laboratories (GCAL) for the samples collected from the Skinner Landfill site in March 2004, was conducted by Earth Tech using the National Functional Guidelines for Organic Data Review, (US EPA, October, 1999) as appropriate. The results were reported by GCAL under SDG 204032408.

| GCAL # | Sample Description |
|-------------|--------------------|
| 20403240801 | SKSWEB1009 |
| 20403240802 | SKSW511009 |
| 20403240803 | SKSW51MS1009 |
| 20403240804 | SKSW51MSD1009 |
| 20403240806 | SKSW521009 |
| 20403240807 | SKSW531009 |

INTRODUCTION

Analyses were performed according to CLP-Organic Analysis Multi-Media, Multi-Concentration OLM04.2 SOW. Results of the sample analyses are reported by the laboratory as either qualified or unqualified. Unqualified results mean that the reported values may be used without reservation. Various data qualifier codes are used by the laboratory to denote specific information regarding the analytical results. The data validation process is intended to evaluate the data on a technical basis. The data package also was subjected to an internal laboratory quality review prior to submission to Earth Tech for data validation.

During the validation process, laboratory-qualified and unqualified data are verified against all available supporting documentation. Based on this evaluation, qualifier codes may be added, deleted or modified by the data user. Final results are therefore, either qualified or unqualified. Validator-qualified results are annotated with the following codes in accordance with the Functional Guidelines:

- U The constituent was analyzed for, but was not detected above the level of the associated analytical reporting limit. The associated value is either the sample quantitation limit or the sample detection limit.
- J The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample.
- UJ The analyte was not detected above the reported sample quantitation limit. However, the reported quantitation limit is approximate and may or may not represent the actual limit of quantitation necessary to accurately and precisely measure the analyte in the sample.

- R The sample results are rejected due to serious deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte cannot be verified.

Details of the semivolatile data validation findings and conclusions are provided in the following sections of this report:

1. Holding Times
2. GC/MS Tuning
3. Calibration
 - A. IC
 - B. CC
4. Blanks
5. System Monitoring Compound Recovery
6. MS/MSD
7. Internal Standards Performance
8. Compound Identification
9. Constituent Quantitation and Reported Detection Limits
10. System Performance
11. Documentation
12. Overall Assessment

1. HOLDING TIMES

All samples were initially extracted within the seven-day technical holding time and the five-day VTSR method holding time. The cooler temperature upon receipt at the laboratory was within the recommended temperature of 4°C +/- 2°C.

2. GC/MS TUNING

The samples were analyzed on a single GC/MS system, identified as MSSV2. Two decafluorotriphenylphosphine (DFTPP) tunes were run representing the shift in which the standards and samples were analyzed. The DFTPP tunes are acceptable.

3. CALIBRATION

A. Initial Calibration

One IC dated 3/19/04 was analyzed in support of the semivolatile sample analyses. Documentation of the IC was present in the data package, and the Relative Response Factor (RRF), as well as percent % RSD values were accurately reported for all target compounds. The criteria employed for technical data review purposes are different than those used in the method. The laboratory must meet a minimum RRF of 0.01; however, for data review purposes, a RRF criterion of "greater than or equal to 0.05" is applied to all semi-volatile compounds. The RRF's and the average RRF were within the acceptance criteria specified in the method for all reported analytes. The %RSD's were within the acceptance criteria specified in the method for all target analytes with the exception of Di-n-butylphthalate (32.0%), Di-n-octylphthalate (30.3%) and Diethylphthalate (41.8%). The lowest point of the calibration curve was dropped for Di-n-butylphthalate, Di-n-octylphthalate, and Diethylphthalate and the %RSD were recalculated. The recalculated %RSD were within the acceptance criteria of less than 30%. Di-n-octylphthalate and Diethylphthalate were not detected in the associated samples therefore data qualification was not required. The detected results for Di-n-butylphthalate were mitigated do to the presence of Di-n-butylphthalate in the associated method blank.

B. Continuing Calibration

One CC dated 4/7/04 was analyzed in support of the semivolatile sample analyses reported in the data submissions. The RRF's for the CC dated 4/7/04 were within the acceptance criteria. The percent difference (%D) between the average RRF's and the CC Response Factors for the CC dated 4/7/04 were within the acceptance criteria with the exception the %D for Di-n-butylphthalate and Di-n-octylphthalate. As per the National Functional Guidelines, if the %D exceeds the acceptance criteria qualify detected results for that analyte with "J" and non-detected results for that analyte with "UJ".

4. BLANKS

One laboratory semivolatile method blank and equipment blank were analyzed with this SDG. The results are summarized below.

Method Blank (0325SBLK)

Di-n-butylphthalate was detected at a concentration of 0.851 ppb in method blank 0322SBLK.

Equipment Blank (SKSWEB1009)

Di-n-butylphthalate was detected at a concentration of 0.904 ppb in the equipment blank collected on 3/22/04. The Di-n-butylphthalate result was mitigated by the presence of Di-n-butylphthalate in the associated extraction blank.

5. SYSTEM MONITORING COMPOUND RECOVERY

All reported semivolatile system monitoring compounds were recovered within acceptable control limits.

6. MATRIX SPIKE/MATRIX SPIKE DUPLICATE (MS/MSD)

Sample SKSW511009 was submitted for MS/MSD analysis. The MS/MSD percent recoveries were within the acceptance criteria with the exception of the 4-Nitorphenol. The %RPD between the MS/MSD are within the acceptance criteria. As per the National Functional Guidelines, no action is taken on MS/MSD data alone.

7. INTERNAL STANDARDS PERFORMANCE

Internal standard areas and retention times were within acceptable limits for the reported semivolatile sample analyses.

8. COMPOUND IDENTIFICATION

All reported semivolatile constituents were correctly identified with supporting chromatograms present in the data package.

9. CONSTITUENT QUANTITATION AND REPORTED DETECTION LIMITS

Constituent quantitations were correctly calculated and reported for semivolatile constituents.

10. SYSTEM PERFORMANCE

The analytical system appears to have been working well at the time of these analyses, based on the evaluation of the raw data submitted for review.

11. DOCUMENTATION

The documentation appeared accurate and in order with the exception of the following: GCAL inadvertently left the Date Extracted off of the CLP Form 1's therefore the data validator inserted the Date Extracted on the CLP Form 1's. GCAL also inadvertently left the "B" qualifier off of the CLP Form 1's for the compound Di-n-butylphthalate therefore the data validator inserted a "B" qualifier in the "Q" column of the CLP Form 1's. The "B" qualifier indicates that the analyte was detected in the associated method blank.

12. OVERALL ASSESSMENT

The results are acceptable with the validator-added qualifiers.

**DATA VALIDATION SUMMARY – SAMPLE DELIVERY GROUP 204032408
VOLATILE ORGANIC**

Validation of the GC/MS volatile organics data, as prepared by Gulf Coast Analytical Laboratories (GCAL) for the samples collected from the Skinner Landfill site in March 2004, was conducted by Earth Tech using the National Functional Guidelines for Organic Data Review, (US EPA, October, 1999), as appropriate. The results were reported by GCAL under SDG 204032408.

| GCAL # | Sample Description |
|-------------|--------------------|
| 20403240801 | SKSWEB1009 |
| 20403240802 | SKSW511009 |
| 20403240803 | SKSW51MS1009 |
| 20403240804 | SKSW51MSD1009 |
| 20403240806 | SKSW521009 |
| 20403240807 | SKSW531009 |

INTRODUCTION

Analyses were performed according to CLP-Organic Analysis Low Concentration OLC02.0 SOW. Results of the sample analyses are reported by the laboratory as either qualified or unqualified. Unqualified results mean that the reported values may be used without reservation. Various qualifier codes are used by the laboratory to denote specific information regarding the analytical results. The data validation process is intended to evaluate the data on a technical basis. The data package also was subjected to an internal laboratory quality review prior to submission to Earth Tech for data validation.

During the validation process, laboratory-qualified and unqualified data are verified against all available supporting documentation. Based on this evaluation, qualifier codes may be added, deleted or modified by the data user. Final results are therefore, either qualified or unqualified. Validator-qualified results are annotated with the following codes in accordance with the Functional Guidelines:

- U The constituent was analyzed for, but was not detected above the level of the associated analytical reporting limit. The associated value is either the sample quantitation limit or the sample detection limit.
- J The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample.
- UJ The analyte was not detected above the reported sample quantitation limit.

However, the reported quantitation limit is approximate and may or may not represent the actual limit of quantitation necessary to accurately and precisely measure the analyte in the sample.

- R The sample results are rejected due to serious deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte cannot be verified.

The volatiles data validation findings and conclusions are provided in the following sections of this report:

1. Holding Times
2. GC/MS Tuning
3. Calibration
 - A. IC
 - B. CC
4. Blanks
5. System Monitoring Compound Recovery
6. MS/MSD
7. Laboratory Control Sample
8. Internal Standards Performance
9. Compound Identification
10. Constituent Quantitation and Reported Detection Limits
11. System Performance
12. Documentation
13. Overall Assessment

1. **HOLDING TIMES**

All samples for Volatile Organic Compounds (VOC) analyses were analyzed within the 14-day technical holding time and the 10-day VTSR method holding time. The cooler temperature upon receipt at the laboratory was within the recommended temperature of 4°C +/- 2°C.

2. GC/MS TUNING

The samples were analyzed on two GC/MS system, identified as MSV0 and MSV2. Two bromofluorobenzene (BFB) tunes were run. The BFB tunes are acceptable.

3. CALIBRATION

A. Initial Calibration

One IC dated 3/23/04 was analyzed on instrument MSV2 in support of the volatile sample analyses reported in the data submissions. Documentation of the IC standards is present in the data package, and RRF's as well as %RSD values were accurately reported. The criteria employed for technical data review purposes are different than those used in the method. The laboratory must meet a minimum RRF of 0.01; however, for data review purposes, a RRF criterion of "greater than or equal to 0.05" is applied to all volatile compounds. The %RSD's were within the acceptance criteria specified in the method for all target analytes. The RRF's and the average RRF for the IC dated 3/23/04 were within the acceptance criteria specified in the method for all target analytes with the exception of Acetone. As per the National Functional Guidelines, if any initial calibration RRF is less than 0.05, qualify positive results that have acceptable mass spectral identification with "J", using professional judgement, and non-detected analytes as unusable (R). It should be noted that the laboratory did meet the minimum RRF of 0.01 for all target compounds.

B. Continuing Calibration

One CC dated 3/23/04 was analyzed on instrument MSV2 in support of the volatile sample analyses reported in the data submissions. The percent difference (%D) between the average RRF's and the CC RF's were within the acceptance criteria for all target analytes. The CC RRF's were within the acceptance criteria specified in the method for all target analytes with the exception of Acetone. The Acetone results were previously qualified under section 3A above.

4. BLANKS

One laboratory volatile method blank, storage blank, and Equipment Blank were analyzed with this SDG. The results are summarized below.

MB154405

1,2-Dichlorobenzene was detected at a concentration of 0.17 ppb in the method blank analyzed on 3/24/04.

Storage Blank (VHBLK01)

No compounds were detected above the MDL in the storage blank analyzed on 3/24/04.

Trip Blank

There was no Trip Blank submitted for this sampling event.

Equipment Blank (SKSWEB1009)

Acetone, Methylene chloride and Toluene were detected at concentrations of 2.8 ppb, 0.47 ppb and 1.1 ppb respectively in the equipment blank collected on 3/22/04.

5. SYSTEM MONITORING COMPOUND RECOVERY

All reported volatile system monitoring compounds were recovered within acceptable control limits for all samples.

6. MATRIX SPIKE/MATRIX SPIKE DUPLICATE

Sample SKSW511009 was submitted for MS/MSD analysis. The MS/MSD percent recoveries and %RPD between the MS/MSD were within the acceptance criteria.

7. LABORATORY CONTROL SAMPLE

A LCS was analyzed in conjunction with this SDG. Recoveries were within the control limit for all constituents.

8. INTERNAL STANDARDS PERFORMANCE

Internal Standard (IS) areas and retention times were within acceptable limits for the reported volatile sample analyses.

9. COMPOUND IDENTIFICATION

All reported VOCs were correctly identified with supporting chromatograms present in the data package.

10. CONSTITUENT QUANTITATION AND REPORTED DETECTION LIMITS

Constituent quantitations were correctly calculated and reported for VOCs with the exception of Ethylbenzene. The Ethylbenzene standard and detected results were originally quantitated using the incorrect quantitation ion (GCAL used 106 instead of 91). GCAL corrected the mistake and re-submitted the corrected pages that were affected in the laboratory report. The overall effect had no impact in the final result for Ethylbenzene.

11. SYSTEM PERFORMANCE

The analytical system appears to have been working well at the time of these analyses, based on the evaluation of the raw data.

12. DOCUMENTATION

The documentation appeared accurate and in order with the exception of the following: GCAL inadvertently transposed the area counts and retention times for IS Chlorobenzene-d5 and 1,4-Dichlorobenzene-d4. The data validator corrected the mistake by drawing arrows to indicate the correct area counts and retention times for IS Chlorobenzene-d5 and 1,4-Dichlorobenzene-d4.

13. OVERALL ASSESSMENT

The results are acceptable with the validator-added qualifiers.

DATA VALIDATION SUMMARY - SAMPLE DELIVERY GROUP 204032408 PESTICIDES

Validation of the Gas Chromatography (GC) pesticides data, as prepared by Gulf Coast Analytical Laboratories (GCAL) for the samples collected from the Skinner Landfill site in March 2004, was conducted by Earth Tech using the National Functional Guidelines for Organic Data Review, (US EPA, October, 1999), as appropriate. The results were reported by GCAL under SDG 204032408.

| GCAL # | Sample Description |
|-------------|--------------------|
| 20403240801 | SKSWEB1009 |
| 20403240802 | SKSW511009 |
| 20403240803 | SKSW51MS1009 |
| 20403240804 | SKSW51MSD1009 |
| 20403240806 | SKSW521009 |
| 20403240807 | SKSW531009 |

INTRODUCTION

Analyses were performed according to CLP-Organic Analysis Multi-Media, Multi-Concentration OLM04.2 SOW. Results of the sample analyses are reported by the laboratory as either qualified or unqualified. Unqualified results mean that the reported values may be used without reservation. Various qualifier codes are used by the laboratory to denote specific information regarding the analytical results.

The data validation process is intended to evaluate the data on a technical basis. The data package also was subjected to an internal laboratory quality review prior to submission to Earth Tech for data validation.

During the validation process, laboratory-qualified and unqualified data are verified against all available supporting documentation. Based on this evaluation, qualifier codes may be added, deleted or modified by the data user. Final results are therefore, either qualified or unqualified. Validator-qualified results are annotated with the following codes in accordance with the Functional Guidelines:

- U The constituent was analyzed for, but was not detected above the level of the associated analytical reporting limit. The associated value is either the sample quantitation limit or the sample detection limit.
- J The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample.
- UJ The analyte was not detected above the reported sample quantitation limit. However, the reported quantitation limit is approximate and may or may not represent the actual limit of quantitation necessary to accurately and precisely measure the analyte in the sample.

- R The sample results are rejected due to serious deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte cannot be verified.

Details of the pesticide data validation findings and conclusions are provided in the following sections of this report:

1. Holding Times
2. Gas Chromatograph/Electronic Capture Detector (GC/ECD) Instrument Performance Check
3. IC
4. Calibration Verification
5. Blanks
6. Surrogate Spikes
7. Matrix Spike/Matrix Spike Duplicate (MS/MSD)
8. Pesticide Cleanup Checks
9. Target Compound Identification
10. Constituent Quantitation and Reported Detection Limits
11. Documentation
12. Overall Assessment

1. HOLDING TIMES

All samples were extracted within the seven-day technical holding time and the five-day VTSR method holding time. The cooler temperature upon receipt at the laboratory was within the recommended temperature of 4°C +/- 2°C.

2. GC/ECD INSTRUMENT PERFORMANCE CHECK

The Performance Evaluation Mixture (PEM) was analyzed at the correct frequency. Absolute retention times were within limits.

The percent resolution between adjacent peaks was within QC limits for the Pesticide Analyte Resolution Check.

The percent resolution between adjacent peaks is within QC limits for the Performance Evaluation Mixtures (PEM). The percent breakdown for both 4,4-DDT and Endrin in each PEM was less than 20.0% for both GC columns. The combined percent breakdown for 4,4-DDT and Endrin in each PEM was less than 30.0% for both GC columns.

3. INITIAL CALIBRATION

Individual standard mixtures A and B were analyzed at the correct frequencies and concentrations. The percent resolution criterion was met for Individual standard mixtures A and B.

The Percent Relative Standard Deviation (%RSD) of the calibration factors for each of the single component pesticides was less than 20%.

The multi-component target compounds were analyzed separately on both columns at a single concentration level. Retention times were determined from a minimum of three peaks.

4. CALIBRATION VERIFICATION

Absolute retention times were within appropriate time retention windows

5. BLANKS

One laboratory method blank and Equipment Blank were analyzed with this SDG. The results are summarized below.

Method Blank 155051

No constituents were detected above the laboratory-reporting limit. This blank corresponds to all samples extracted on 3/26/04.

Equipment Blank SKSWEB1009

No constituents were detected above the laboratory-reporting limit in the equipment blank collected on 3/22/04.

6. SURROGATE SPIKES

Decachlorobiphenyl (DCB) and tetrachloro-m-xylene (TCX) surrogate spike recoveries were within the acceptance criteria for all samples.

7. MATRIX SPIKE/MATRIX SPIKE DUPLICATE

Sample SKSW511009 was submitted for MS/MSD analysis. The MS/MSD percent recoveries were within the acceptance criteria. The %RPD between the MS/MSD are within the acceptance criteria.

8. PESTICIDE CLEANUP CHECKS

Recoveries of all pesticides and surrogates were within 80-120% for the lot of Florisil cartridges utilized for pesticide cleanup with the exception of 4,4'-DDT (130%). There were no target compounds detected in the associated samples therefore no action was taken.

9. TARGET COMPOUND IDENTIFICATION

All reported pesticide data were correctly identified with supporting chromatograms present in the data package.

10. CONSTITUENT QUANTITATION AND REPORTED DETECTION LIMITS

Constituent quantitations were correctly calculated and reported for pesticide constituents.

11. DOCUMENTATION

The documentation appeared accurate and in order with the exception of the following: GCAL inadvertently left the Date Extracted off of the CLP Form 1's therefore the data validator inserted the Date Extracted on the CLP Form 1's.

12. OVERALL ASSESSMENT

The results are acceptable with the validator-added qualifiers.

REFERENCES

US EPA, 1994. *National Functional Guidelines for Inorganic Data Review.*

US EPA, 1999. *National Functional Guidelines for Organic Data Review.*

ANALYTICAL RESULTS

PERFORMED BY

GULF COAST ANALYTICAL LABORATORIES, INC.

Report Date 04/16/2004

GCAL Report 204032408

RESUBMITTED

Deliver To Earth Tech
200 Vine Street
Wilder, KY 41076
859-442-2300

Attn Pat Higgins

Customer Earth Tech

Project Skinner Landfill

CASE NARRATIVE

Client: Earth Tech **Report:** 204032408

Gulf Coast Analytical Laboratories received and analyzed the sample(s) listed on the sample cross-reference page of this report. Receipt of the sample(s) is documented by the attached chain of custody. This applies only to the sample(s) listed in this report. No sample integrity or quality control exceptions were identified unless noted below.

In the ILM04.1 CLP metals, Forms I, II, III, V, VI, VII, IX, X, and XIV were resubmitted 07/28/2005 to report TAL metals.

SEMI-VOLATILES MASS SPECTROMETRY

The MS/MSD recoveries for 4-Nitrophenol are above the control limits.

SEMI-VOLATILES GAS CHROMATOGRAPHY

In the Florisil check analysis, the recovery for DDT was above recovery limit, however DDT was not detected in the associated samples.

METALS

Barium, Chromium, Iron, and Nickel are flagged as estimated for samples associated with prep batch 271292 due to the fact that the percent difference between the original sample result and the serial dilution result is greater than 10. A chemical or physical interference is suspected. In the ILM04.1 - CLP Metals analysis, the MS recovery was outside the control limits for Antimony, Arsenic, Selenium, and Zinc. The LCS recovery was within control limits. This indicates the analysis is in control and the sample is affected by matrix interference. A post-digestion spike was performed on the QC sample for this batch with a recovery of 109% for Antimony, 131% for Arsenic, 64% for Selenium, and 139% for Zinc. The MS recovery is not applicable for Aluminum, Manganese, Iron, and Lead because the sample concentration is greater than four times the spike concentration. The Sample/Duplicate RPD for Aluminum, Arsenic, Barium, Calcium, Chromium, Cobalt, Copper, Iron, Lead, Magnesium, Manganese, Nickel, Potassium, Vanadium, and Zinc was outside the control limits. The heterogeneous nature of the QC sample is believed to be responsible for this.

In the ILM04.1 - CLP Metals analysis for prep batch 271294, the MS recovery was outside the control limits for Arsenic and Selenium. The LCS recovery was within control limits. This indicates the analysis is in control and the sample is affected by matrix interference. A post-digestion spike was performed on the QC sample for this batch with a recovery of 106% for Arsenic and 94% for Selenium. The Sample/Duplicate RPD for Vanadium is not applicable because the sample and/or duplicate concentration are less than five times the reporting limit.

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RESUBMITTED

U.S. EPA - CLP
1
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKSWEB1009

Lab Name: GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204032408
Matrix: (soil / water) Water Lab Sample ID: 20403240801
Level: (low / med) _____ Date Received: 03/24/04
% Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M | |
|-----------|-----------|---------------|---|-----|----|----|
| 7429-90-5 | Aluminum | 25.8 | U | *E | P | |
| 7440-36-0 | Antimony | 3.7 | U | N | P | UJ |
| 7440-38-2 | Arsenic | 2.9 | U | N,* | P | |
| 7440-39-3 | Barium | 0.3 | B | *E | P | J |
| 7440-41-7 | Beryllium | 0.1 | U | / | P | |
| 7440-43-9 | Cadmium | 0.2 | U | / | P | |
| 7440-70-2 | Calcium | 34.9 | B | * | P | J |
| 7440-47-3 | Chromium | 1.0 | B | *E | P | J |
| 7440-48-4 | Cobalt | 0.4 | B | / | P | |
| 7440-50-8 | Copper | 6.5 | B | * | P | J |
| 7439-89-5 | Iron | 37.8 | B | *E | P | J |
| 7439-92-1 | Lead | 1.5 | U | * | P | |
| 7439-95-4 | Magnesium | 36.7 | U | *E | P | |
| 7439-96-5 | Manganese | 0.4 | B | * | P | J |
| 7439-97-6 | Mercury | 0.1 | U | / | AV | |
| 7440-02-0 | Nickel | 1.4 | B | *E | P | J |
| 7440-09-7 | Potassium | 42.1 | U | /E | P | J |
| 7782-49-2 | Selenium | 4.4 | U | N | P | R |
| 7440-22-4 | Silver | 1.2 | B | | P | |
| 7440-23-5 | Sodium | 45.4 | U | | P | |
| 7440-28-0 | Thallium | 2.6 | U | | P | |
| 7440-62-2 | Vanadium | 0.8 | U | E | P | |
| 7440-66-6 | Zinc | 9.7 | B | N,* | P | J |
| 57-12-5 | Cyanide | 0.5 | U | | AS | |

3/24/05
msr

Color Before: COLORLESS Clarity Before: CLEAR Texture: _____
Color After: COLORLESS Clarity After: CLEAR Artifacts: _____
Comments: _____

U.S. EPA - CLP
1
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKSW511009

50

Lab Name: GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204032408
Matrix: (soil / water) Water Lab Sample ID: 20403240802
Level: (low / med) _____ Date Received: 03/24/04
% Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|-----|----|
| 7429-90-5 | Aluminum | 62300 | | *,E | P |
| 7440-36-0 | Antimony | 7.3 | B | N | P |
| 7440-38-2 | Arsenic | 50.7 | | N,* | P |
| 7440-39-3 | Barium | 499 | | *,E | P |
| 7440-41-7 | Beryllium | 4.9 | B | / | P |
| 7440-43-9 | Cadmium | 5.0 | | / | P |
| 7440-70-2 | Calcium | 427000 | | * | P |
| 7440-47-3 | Chromium | 72.6 | | *,E | P |
| 7440-48-4 | Cobalt | 59.7 | | / | P |
| 7440-50-8 | Copper | 131 | | * | P |
| 7439-89-6 | Iron | 124000 | | *,E | P |
| 7439-92-1 | Lead | 122 | | * | P |
| 7439-95-4 | Magnesium | 80300 | | *,E | P |
| 7439-96-5 | Manganese | 5690 | | * | P |
| 7439-97-6 | Mercury | 0.1 | B | / | AV |
| 7440-02-0 | Nickel | 116 | | *,E | P |
| 7440-09-7 | Potassium | 12200 | | /,E | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-23-5 | Sodium | 60200 | | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-62-2 | Vanadium | 105 | | E | P |
| 7440-66-6 | Zinc | 490 | | N,* | P |
| 57-12-5 | Cyanide | 0.8 | B | | AS |

8/29/05
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Color Before: DK.BROWN Clarity Before: CLOUDY Texture: _____
Color After: LT.BROWN Clarity After: CLEAR Artifacts: _____
Comments: _____

EPA SAMPLE NO.

SKSW51MS1009

Lab Name: GCAL Contract:

Lab Code: LA024 Case No.: SAS No.: SDG No.: 204032408

Matrix: (soil / water) Water Lab Sample ID: 20403240803

Level: (low / med) _____ Date Received: 03/24/04

% Solids:

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|-----|----|
| 7429-90-5 | Aluminum | 82700 | | *,E | P |
| 7440-36-0 | Antimony | 41.9 | B | N | P |
| 7440-38-2 | Arsenic | 111 | | N,* | P |
| 7440-39-3 | Barium | 2580 | | *,E | P |
| 7440-41-7 | Beryllium | 56.9 | | / | P |
| 7440-43-9 | Cadmium | 51.6 | | / | P |
| 7440-70-2 | Calcium | 587000 | | * | P |
| 7440-47-3 | Chromium | 284 | | *,E | P |
| 7440-48-4 | Cobalt | 545 | | / | P |
| 7440-50-8 | Copper | 436 | | * | P |
| 7439-89-6 | Iron | 168000 | | *,E | P |
| 7439-92-1 | Lead | 183 | | * | P |
| 7439-95-4 | Magnesium | 101000 | | *,E | P |
| 7439-96-5 | Manganese | 8350 | | * | P |
| 7439-97-6 | Mercury | 5.8 | | / | AV |
| 7440-02-0 | Nickel | 619 | | *,E | P |
| 7440-09-7 | Potassium | 13900 | | /,E | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 46.3 | | | P |
| 7440-23-5 | Sodium | 58900 | | | P |
| 7440-28-0 | Thallium | 56.7 | | | P |
| 7440-62-2 | Vanadium | 614 | | E | P |
| 7440-66-6 | Zinc | 1130 | | N,* | P |
| 57-12-5 | Cyanide | 87.8 | | | AS |

[illegible]

5/25/17
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Color Before: DK.BROWN

Clarity Before: CLOUDY

Texture:

Color After: LT.BROWN

Clarity After: CLEAR

Artifacts:

Comments:

U.S. EPA - CLP
1
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKSW51DUP1009
50

Lab Name: GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204032408
Matrix: (soil / water) Water Lab Sample ID: 20403240805
Level: (low / med) _____ Date Received: 03/24/04
% Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|-----|----|
| 7429-90-5 | Aluminum | 88100 | | *E | P |
| 7440-36-0 | Antimony | 6.7 | B | N | P |
| 7440-38-2 | Arsenic | 67.0 | | N,* | P |
| 7440-39-3 | Barium | 714 | | *E | P |
| 7440-41-7 | Beryllium | 6.7 | | / | P |
| 7440-43-9 | Cadmium | 7.2 | | / | P |
| 7440-70-2 | Calcium | 603000 | | * | P |
| 7440-47-3 | Chromium | 102 | | *E | P |
| 7440-48-4 | Cobalt | 83.4 | | / | P |
| 7440-50-8 | Copper | 174 | | * | P |
| 7439-89-6 | Iron | 173000 | | *E | P |
| 7439-92-1 | Lead | 169 | | * | P |
| 7439-95-4 | Magnesium | 102000 | | *E | P |
| 7439-96-5 | Manganese | 7950 | | * | P |
| 7439-97-6 | Mercury | 0.3 | | / | AV |
| 7440-02-0 | Nickel | 162 | | *E | P |
| 7440-09-7 | Potassium | 17000 | | /E | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-23-5 | Sodium | 60100 | | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-62-2 | Vanadium | 145 | | E | P |
| 7440-66-6 | Zinc | 679 | | N,* | P |
| 57-12-5 | Cyanide | 1.1 | B | | AS |

8/29/05
msl

Color Before: DK.BROWN Clarity Before: CLOUDY Texture: _____
Color After: LT.BROWN Clarity After: CLEAR Artifacts: _____
Comments: _____

U.S. EPA - CLP
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INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKSW521009

51

Lab Name: GCAL Contract: _____
 Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204032408
 Matrix: (soil / water) Water Lab Sample ID: 20403240806
 Level: (low / med) _____ Date Received: 03/24/04
 % Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|-----|----|
| 7429-90-5 | Aluminum | 9250 | | *,E | P |
| 7440-36-0 | Antimony | 3.7 | U | N | P |
| 7440-38-2 | Arsenic | 11.1 | | N,* | P |
| 7440-39-3 | Barium | 112 | B | *,E | P |
| 7440-41-7 | Beryllium | 0.7 | B | / | P |
| 7440-43-9 | Cadmium | 0.5 | B | / | P |
| 7440-70-2 | Calcium | 153000 | | * | P |
| 7440-47-3 | Chromium | 12.7 | | *,E | P |
| 7440-48-4 | Cobalt | 7.8 | B | / | P |
| 7440-50-8 | Copper | 22.0 | B | * | P |
| 7439-89-6 | Iron | 17800 | | *,E | P |
| 7439-92-1 | Lead | 17.3 | | * | P |
| 7439-95-4 | Magnesium | 38900 | | *,E | P |
| 7439-96-5 | Manganese | 685 | | * | P |
| 7439-97-6 | Mercury | 0.1 | U | / | AV |
| 7440-02-0 | Nickel | 16.4 | B | *,E | P |
| 7440-09-7 | Potassium | 4470 | B | /,E | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-23-5 | Sodium | 61800 | | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-62-2 | Vanadium | 18.7 | B | E | P |
| 7440-66-6 | Zinc | 52.9 | | N,* | P |
| 57-12-5 | Cyanide | 0.6 | B | | AS |

8/29/05
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Color Before: LT.BROWN Clarity Before: CLEAR Texture: _____
 Color After: LT.BROWN Clarity After: CLEAR Artifacts: _____
 Comments: _____

U.S. EPA - CLP
1
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKSW531009
52

Lab Name: GCAL

Contract: _____

Lab Code: LA024

Case No.: _____

SAS No.: _____

SDG No.: 204032408

Matrix: (soil / water) Water

Lab Sample ID: 20403240807

Level: (low / med) _____

Date Received: 03/24/04

% Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|----|----|
| 7429-90-5 | Aluminum | 26.8 | B | *E | P |
| 7440-36-0 | Antimony | 3.7 | U | N | P |
| 7440-38-2 | Arsenic | 2.9 | U | N* | P |
| 7440-39-3 | Barium | 41.7 | B | *E | P |
| 7440-41-7 | Beryllium | 0.1 | U | / | P |
| 7440-43-9 | Cadmium | 0.2 | U | / | P |
| 7440-70-2 | Calcium | 108000 | | * | P |
| 7440-47-3 | Chromium | 1.7 | B | *E | P |
| 7440-48-4 | Cobalt | 0.5 | B | / | P |
| 7440-50-8 | Copper | 3.2 | B | * | P |
| 7439-89-6 | Iron | 67.8 | B | *E | P |
| 7439-92-1 | Lead | 1.5 | U | * | P |
| 7439-95-4 | Magnesium | 31100 | | *E | P |
| 7439-96-5 | Manganese | 3.2 | B | * | P |
| 7439-97-6 | Mercury | 0.1 | U | / | AV |
| 7440-02-0 | Nickel | 1.1 | B | *E | P |
| 7440-09-7 | Potassium | 1900 | B | /E | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.6 | B | | P |
| 7440-23-5 | Sodium | 61200 | | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-62-2 | Vanadium | 2.2 | B | E | P |
| 7440-66-6 | Zinc | 0.6 | U | N* | P |
| 57-12-5 | Cyanide | 0.5 | U | | AS |

8/29/05
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Color Before: COLORLESS

Clarity Before: CLEAR

Texture: _____

Color After: COLORLESS

Clarity After: CLEAR

Artifacts: _____

Comments:

U.S. EPA - CLP
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INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKSWEB1009 (DISS)

Lab Name: GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204032408
Matrix: (soil / water) Water Lab Sample ID: 20403240808
Level: (low / med) _____ Date Received: 03/24/04
% Solids: _____
Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum | 25.8 | U | | P |
| 7440-36-0 | Antimony | 3.7 | U | | P |
| 7440-38-2 | Arsenic | 2.9 | U | N | P |
| 7440-39-3 | Barium | 0.4 | B | | P |
| 7440-41-7 | Beryllium | 0.1 | U | | P |
| 7440-43-9 | Cadmium | 0.2 | U | | P |
| 7440-70-2 | Calcium | 60.8 | B | | P |
| 7440-47-3 | Chromium | 0.8 | U | | P |
| 7440-48-4 | Cobalt | 0.4 | U | | P |
| 7440-50-8 | Copper | 4.2 | B | | P |
| 7439-89-6 | Iron | 14.1 | U | | P |
| 7439-92-1 | Lead | 1.5 | U | | P |
| 7439-95-4 | Magnesium | 36.7 | U | | P |
| 7439-96-5 | Manganese | 0.2 | U | E | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 0.7 | U | | P |
| 7440-09-7 | Potassium | 42.1 | U | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-23-5 | Sodium | 308 | B | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-62-2 | Vanadium | 0.8 | U | | P |
| 7440-66-6 | Zinc | 5.6 | B | | P |

3/29/05
ms

Color Before: COLORLESS Clarity Before: CLEAR Texture: _____
Color After: COLORLESS Clarity After: CLEAR Artifacts: _____
Comments:

U.S. EPA - CLP
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INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKSW511009 (DISS)
50

Lab Name: GCAL Contract: _____
 Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204032408
 Matrix: (soil / water) Water Lab Sample ID: 20403240809
 Level: (low / med) _____ Date Received: 03/24/04
 % Solids: _____
 Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum | 25.8 | U | | P |
| 7440-36-0 | Antimony | 3.7 | U | | P |
| 7440-38-2 | Arsenic | 2.9 | U | N | P |
| 7440-39-3 | Barium | 35.8 | B | | P |
| 7440-41-7 | Beryllium | 0.1 | U | | P |
| 7440-43-9 | Cadmium | 0.2 | U | | P |
| 7440-70-2 | Calcium | 103000 | | | P |
| 7440-47-3 | Chromium | 1.6 | B | | P |
| 7440-48-4 | Cobalt | 0.4 | B | | P |
| 7440-50-8 | Copper | 4.3 | B | | P |
| 7439-89-6 | Iron | 14.1 | U | | P |
| 7439-92-1 | Lead | 1.5 | U | | P |
| 7439-95-4 | Magnesium | 29700 | | | P |
| 7439-96-5 | Manganese | 30.0 | | E | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 1.0 | B | | P |
| 7440-09-7 | Potassium | 2980 | B | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-23-5 | Sodium | 59200 | | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-62-2 | Vanadium | 2.2 | B | | P |
| 7440-66-6 | Zinc | 0.6 | U | | P |

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Color Before: COLORLESS Clarity Before: CLEAR Texture: _____
 Color After: COLORLESS Clarity After: CLEAR Artifacts: _____
 Comments: _____

U.S. EPA - CLP
1
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKSW51MS1009 (DISS)
50

Lab Name: GCAL

Contract: _____

Lab Code: LA024

Case No.: _____

SAS No.: _____

SDG No.: 204032408

Matrix: (soil / water) Water

Lab Sample ID: 20403240810

Level: (low / med) _____

Date Received: 03/24/04

% Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum | 2190 | | | P |
| 7440-36-0 | Antimony | 124 | | | P |
| 7440-38-2 | Arsenic | 51.2 | | N | P |
| 7440-39-3 | Barium | 2140 | | | P |
| 7440-41-7 | Beryllium | 55.8 | | | P |
| 7440-43-9 | Cadmium | 52.5 | | | P |
| 7440-70-2 | Calcium | 105000 | | | P |
| 7440-47-3 | Chromium | 212 | | | P |
| 7440-48-4 | Cobalt | 515 | | | P |
| 7440-50-8 | Copper | 291 | | | P |
| 7439-89-6 | Iron | 1110 | | | P |
| 7439-92-1 | Lead | 19.0 | | | P |
| 7439-95-4 | Magnesium | 30300 | | | P |
| 7439-96-5 | Manganese | 562 | | E | P |
| 7439-97-6 | Mercury | 5.8 | | | AV |
| 7440-02-0 | Nickel | 519 | | | P |
| 7440-09-7 | Potassium | 2790 | B | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 55.0 | | | P |
| 7440-23-5 | Sodium | 59800 | | | P |
| 7440-28-0 | Thallium | 50.2 | | | P |
| 7440-62-2 | Vanadium | 537 | | | P |
| 7440-66-6 | Zinc | 504 | | | P |

Color Before: COLORLESS

Clarity Before: CLEAR

Texture: _____

Color After: COLORLESS

Clarity After: CLEAR

Artifacts: _____

Comments:

U.S. EPA - CLP
1
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKSW5150
DUP1009 (DISS)

Lab Name: GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204032408
Matrix: (soil / water) Water Lab Sample ID: 20403240811
Level: (low / med) _____ Date Received: 03/24/04
% Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum | 25.8 | U | | P |
| 7440-36-0 | Antimony | 3.7 | U | | P |
| 7440-38-2 | Arsenic | 2.9 | U | N | P |
| 7440-39-3 | Barium | 33.0 | B | | P |
| 7440-41-7 | Beryllium | 0.1 | U | | P |
| 7440-43-9 | Cadmium | 0.2 | U | | P |
| 7440-70-2 | Calcium | 101000 | | | P |
| 7440-47-3 | Chromium | 1.4 | B | | P |
| 7440-48-4 | Cobalt | 0.4 | B | | P |
| 7440-50-8 | Copper | 1.6 | B | | P |
| 7439-89-6 | Iron | 14.1 | U | | P |
| 7439-92-1 | Lead | 1.5 | U | | P |
| 7439-95-4 | Magnesium | 29100 | | | P |
| 7439-96-5 | Manganese | 29.7 | | E | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 1.0 | B | | P |
| 7440-09-7 | Potassium | 2670 | B | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-23-5 | Sodium | 57300 | | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-62-2 | Vanadium | 1.7 | B | | P |
| 7440-66-6 | Zinc | 0.6 | U | | P |

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Color Before: COLORLESS Clarity Before: CLEAR Texture: _____
Color After: COLORLESS Clarity After: CLEAR Artifacts: _____
Comments: _____

U.S. EPA - CLP
1
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKSW521009 (DISS)
51

Lab Name: GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204032408
Matrix: (soil / water) Water Lab Sample ID: 20403240812
Level: (low / med) _____ Date Received: 03/24/04
% Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum | 25.8 | U | | P |
| 7440-36-0 | Antimony | 3.7 | U | | P |
| 7440-38-2 | Arsenic | 2.9 | U | N | P |
| 7440-39-3 | Barium | 36.7 | B | | P |
| 7440-41-7 | Beryllium | 0.1 | U | | P |
| 7440-43-9 | Cadmium | 0.2 | U | | P |
| 7440-70-2 | Calcium | 106000 | | | P |
| 7440-47-3 | Chromium | 1.7 | B | | P |
| 7440-48-4 | Cobalt | 0.4 | U | | P |
| 7440-50-8 | Copper | 3.0 | B | | P |
| 7439-89-6 | Iron | 14.1 | U | | P |
| 7439-92-1 | Lead | 1.5 | U | | P |
| 7439-95-4 | Magnesium | 31500 | | | P |
| 7439-96-5 | Manganese | 29.9 | | E | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 2.4 | B | | P |
| 7440-09-7 | Potassium | 2160 | B | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-23-5 | Sodium | 60200 | | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-62-2 | Vanadium | 2.4 | B | | P |
| 7440-66-6 | Zinc | 3.4 | B | | P |

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Color Before: COLORLESS Clarity Before: CLEAR Texture: _____
Color After: COLORLESS Clarity After: CLEAR Artifacts: _____
Comments: _____

U.S. EPA - CLP
1
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKSW581009 (DISS)
52

Lab Name: GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204032408
Matrix: (soil / water) Water Lab Sample ID: 20403240813
Level: (low / med) _____ Date Received: 03/24/04
% Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum | 25.8 | U | | P |
| 7440-36-0 | Antimony | 3.7 | U | | P |
| 7440-38-2 | Arsenic | 2.9 | U | N | P |
| 7440-39-3 | Barium | 40.7 | B | | P |
| 7440-41-7 | Beryllium | 0.1 | U | | P |
| 7440-43-9 | Cadmium | 0.2 | U | | P |
| 7440-70-2 | Calcium | 107000 | | | P |
| 7440-47-3 | Chromium | 1.3 | B | | P |
| 7440-48-4 | Cobalt | 0.4 | U | | P |
| 7440-50-8 | Copper | 2.0 | B | | P |
| 7439-89-6 | Iron | 14.1 | U | | P |
| 7439-92-1 | Lead | 1.5 | U | | P |
| 7439-95-4 | Magnesium | 30700 | | | P |
| 7439-96-5 | Manganese | 1.5 | B | E | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 0.7 | U | | P |
| 7440-09-7 | Potassium | 1900 | B | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-23-5 | Sodium | 61400 | | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-62-2 | Vanadium | 1.9 | B | | P |
| 7440-66-6 | Zinc | 4.5 | B | | P |

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Color Before: COLORLESS Clarity Before: CLEAR Texture: _____
Color After: COLORLESS Clarity After: CLEAR Artifacts: _____
Comments: _____

INITIAL AND CONTINUING CALIBRATION VERIFICATION

Lab Name: GCAL Contract: _____
 Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204032408
 Calibration Source: 106-61-2 CPI Instrument ID: ICP5 ICAL ID: 2
 Date Analyzed: 03/29/04 Time: 1215

CRDL STANDARD

| Analyte | True | Found | CAL %R | Units | Method | Type |
|-----------|------|-------|--------|-------|----------------------|------|
| Antimony | 120 | 133 | 111 | ug/L | ILM04.1 - CLP Metals | P |
| Arsenic | 20.0 | 22.2 | 111 | ug/L | ILM04.1 - CLP Metals | P |
| Beryllium | 10.0 | 9.60 | 96 | ug/L | ILM04.1 - CLP Metals | P |
| Cadmium | 10.0 | 9.70 | 97 | ug/L | ILM04.1 - CLP Metals | P |
| Chromium | 20.0 | 19.8 | 99 | ug/L | ILM04.1 - CLP Metals | P |
| Cobalt | 100 | 94.7 | 95 | ug/L | ILM04.1 - CLP Metals | P |
| Copper | 50.0 | 40.0 | 80 | ug/L | ILM04.1 - CLP Metals | P |
| Lead | 6.00 | 5.10 | 84 | ug/L | ILM04.1 - CLP Metals | P |
| Manganese | 30.0 | 27.6 | 92 | ug/L | ILM04.1 - CLP Metals | P |
| Nickel | 80.0 | 76.7 | 96 | ug/L | ILM04.1 - CLP Metals | P |
| Selenium | 10.0 | 10.3 | 103 | ug/L | ILM04.1 - CLP Metals | P |
| Silver | 20.0 | 19.9 | 100 | ug/L | ILM04.1 - CLP Metals | P |
| Thallium | 20.0 | 22.4 | 112 | ug/L | ILM04.1 - CLP Metals | P |
| Vanadium | 100 | 96.4 | 96 | ug/L | ILM04.1 - CLP Metals | P |
| Zinc | 40.0 | 35.5 | 89 | ug/L | ILM04.1 - CLP Metals | P |

INITIAL AND CONTINUING CALIBRATION VERIFICATION

Lab Name: GCAL

Contract: _____

Lab Code: LA024 Case No.: _____

SAS No.: _____ SDG No.: 204032408

Calibration Source: 106-61-2 CPI

Instrument ID: ICP5 ICAL ID: 2

Date Analyzed: 03/29/04 Time: 1506

CRDL STANDARD

| Analyte | True | Found | CAL %R | Units | Method | Type |
|-----------|------|-------|--------|-------|----------------------|------|
| Antimony | 120 | 130 | 108 | ug/L | ILM04.1 - CLP Metals | P |
| Arsenic | 20.0 | 23.2 | 116 | ug/L | ILM04.1 - CLP Metals | P |
| Beryllium | 10.0 | 9.70 | 97 | ug/L | ILM04.1 - CLP Metals | P |
| Cadmium | 10.0 | 9.60 | 96 | ug/L | ILM04.1 - CLP Metals | P |
| Chromium | 20.0 | 19.9 | 100 | ug/L | ILM04.1 - CLP Metals | P |
| Cobalt | 100 | 94.3 | 94 | ug/L | ILM04.1 - CLP Metals | P |
| Copper | 50.0 | 38.5 | 77 | ug/L | ILM04.1 - CLP Metals | P |
| Lead | 6.00 | 5.40 | 90 | ug/L | ILM04.1 - CLP Metals | P |
| Manganese | 30.0 | 27.9 | 93 | ug/L | ILM04.1 - CLP Metals | P |
| Nickel | 80.0 | 76.8 | 96 | ug/L | ILM04.1 - CLP Metals | P |
| Selenium | 10.0 | 10.0 | 100 | ug/L | ILM04.1 - CLP Metals | P |
| Silver | 20.0 | 21.0 | 105 | ug/L | ILM04.1 - CLP Metals | P |
| Thallium | 20.0 | 19.6 | 98 | ug/L | ILM04.1 - CLP Metals | P |
| Vanadium | 100 | 98.2 | 98 | ug/L | ILM04.1 - CLP Metals | P |
| Zinc | 40.0 | 35.0 | 87 | ug/L | ILM04.1 - CLP Metals | P |

INITIAL AND CONTINUING CALIBRATION VERIFICATION

Lab Name: GCAL Contract: _____
 Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204032408
 Calibration Source: 106-61-2 CPI Instrument ID: ICP5 ICAL ID: 2
 Date Analyzed: 03/29/04 Time: 1605

CRDL STANDARD

| Analyte | True | Found | CAL %R | Units | Method | Type |
|-----------|------|-------|--------|-------|----------------------|------|
| Antimony | 120 | 130 | 109 | ug/L | ILM04.1 - CLP Metals | P |
| Arsenic | 20.0 | 20.4 | 102 | ug/L | ILM04.1 - CLP Metals | P |
| Beryllium | 10.0 | 9.60 | 96 | ug/L | ILM04.1 - CLP Metals | P |
| Cadmium | 10.0 | 9.70 | 97 | ug/L | ILM04.1 - CLP Metals | P |
| Chromium | 20.0 | 20.2 | 101 | ug/L | ILM04.1 - CLP Metals | P |
| Cobalt | 100 | 95.4 | 95 | ug/L | ILM04.1 - CLP Metals | P |
| Copper | 50.0 | 39.1 | 78 | ug/L | ILM04.1 - CLP Metals | P |
| Lead | 6.00 | 5.00 | 83 | ug/L | ILM04.1 - CLP Metals | P |
| Manganese | 30.0 | 27.8 | 93 | ug/L | ILM04.1 - CLP Metals | P |
| Nickel | 80.0 | 77.6 | 97 | ug/L | ILM04.1 - CLP Metals | P |
| Selenium | 10.0 | 9.10 | 91 | ug/L | ILM04.1 - CLP Metals | P |
| Silver | 20.0 | 20.9 | 104 | ug/L | ILM04.1 - CLP Metals | P |
| Thallium | 20.0 | 20.4 | 102 | ug/L | ILM04.1 - CLP Metals | P |
| Vanadium | 100 | 97.2 | 97 | ug/L | ILM04.1 - CLP Metals | P |
| Zinc | 40.0 | 35.3 | 88 | ug/L | ILM04.1 - CLP Metals | P |

INITIAL AND CONTINUING CALIBRATION VERIFICATION

Lab Name: GCAL Contract: _____
 Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204032408
 Calibration Source: 106-58-5 EXAXOL Instrument ID: FIMS1 ICAL ID: 1
 Date Analyzed: 03/31/04 Time: 1540

CRDL STANDARD

| Analyte | True | Found | CAL %R | Units | Method | Type |
|---------|-------|-------|--------|-------|----------------|------|
| Mercury | 0.200 | 0.200 | 106 | ug/L | ILM04.1 CLP HG | AV |

U.S. EPA - CLP

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BLANKS

Lab Name: GCAL

Contract:

Lab Code: LA024

Case No.:

SAS No.:

SDG No.: 204032408

Preparation Blank Matrix: (soil / water) Water

Preparation Blank Concentration Units: (ug/L / mg/kg) ug/L

| Analyte | Initial Calib. Blank (ug/L) | | Continuing Calibration Blank (ug/L) | | | | | | Preparation Blank | | M |
|-----------|-----------------------------|---|-------------------------------------|---|------|---|------|---|-------------------|---|----|
| | | C | 1 | C | 2 | C | 3 | C | | C | |
| Aluminum | 25.8 | U | 25.8 | U | 25.8 | U | 25.8 | U | 25.800 | U | P |
| Antimony | 8.8 | B | 4.3 | B | 3.7 | U | 5.5 | B | 3.700 | U | P |
| Arsenic | 2.9 | U | 2.9 | U | 2.9 | U | 2.9 | U | 2.900 | U | P |
| Barium | 0.3 | U | 0.3 | U | 0.3 | U | 0.3 | U | 0.300 | U | P |
| Beryllium | 0.1 | U | 0.1 | U | 0.1 | U | 0.1 | U | 0.100 | U | P |
| Cadmium | 1.3 | B | 1.1 | B | 0.2 | U | 1.6 | B | 0.200 | U | P |
| Calcium | 7.5 | U | 7.5 | U | 7.5 | U | 7.5 | U | 7.500 | U | P |
| Chromium | 0.8 | U | 0.8 | U | 0.8 | U | 0.8 | U | 1.199 | B | P |
| Cobalt | 0.4 | U | 0.4 | U | 0.4 | U | 0.5 | B | 0.822 | B | P |
| Copper | 19.9 | B | 20.5 | B | 5.2 | B | 9.9 | B | 17.654 | B | P |
| Iron | 14.1 | U | 14.1 | U | 14.1 | U | 14.1 | U | 14.100 | U | P |
| Lead | 1.5 | B | 1.5 | U | 1.5 | U | 1.5 | U | 1.500 | U | P |
| Magnesium | 36.7 | U | 36.7 | U | 36.7 | U | 45.5 | B | 36.700 | U | P |
| Manganese | -1 | B | -1.5 | B | 0.2 | U | -1.7 | B | -1.587 | B | P |
| Mercury | 0.1 | U | 0.1 | U | 0.1 | U | 0.1 | U | 0.100 | U | AV |
| Nickel | 0.7 | U | 0.7 | U | 0.7 | U | 0.7 | U | 1.794 | B | P |
| Potassium | 42.1 | U | 42.1 | U | 42.1 | U | 42.1 | U | 42.100 | U | P |
| Selenium | 4.4 | U | 4.4 | U | 4.4 | U | 4.4 | U | 4.400 | U | P |
| Silver | 0.4 | U | 0.4 | B | 0.7 | B | 0.8 | B | 1.381 | B | P |
| Sodium | 45.4 | U | 45.4 | U | 45.4 | U | 45.4 | U | 45.400 | U | P |
| Thallium | 2.6 | U | 2.6 | U | 2.6 | U | 2.6 | U | 2.600 | U | P |
| Vanadium | 0.8 | U | 0.8 | U | 0.8 | U | 0.8 | U | 0.800 | U | P |
| Zinc | 2.6 | B | 3.6 | B | 1.0 | B | 1.2 | B | 7.236 | B | P |
| Cyanide | 0.5 | U | 0.5 | U | 0.5 | U | 0.5 | U | 0.500 | U | AS |

U.S. EPA - CLP

3

BLANKS

Lab Name: GCAL

Contract: _____

Lab Code: LA024 Case No.: _____SAS No.: _____ SDG No.: 204032408Preparation Blank Matrix: (soil / water) WaterPreparation Blank Concentration Units: (ug/L / mg/kg) ug/L

| Analyte | Initial Calib. Blank (ug/L) | Continuing Calibration Blank (ug/L) | | | | | | Prepa- ration Blank C | M |
|-----------|--------------------------------------|-------------------------------------|---|---|---|---|---|--------------------------------|----|
| | | 4 | C | 5 | C | 6 | C | | |
| Aluminum | | 25.8 | U | | | | | 25.800 | P |
| Antimony | | 5.2 | B | | | | | 3.700 | P |
| Arsenic | | 2.9 | U | | | | | 2.900 | P |
| Barium | | 0.3 | U | | | | | 0.300 | P |
| Beryllium | | 0.1 | U | | | | | 0.100 | P |
| Cadmium | | 1.5 | B | | | | | 0.200 | P |
| Calcium | | 7.5 | U | | | | | 7.500 | P |
| Chromium | | 0.8 | B | | | | | 0.800 | P |
| Cobalt | | 0.4 | U | | | | | 0.475 | P |
| Copper | | 13.5 | B | | | | | 5.096 | P |
| Iron | | -16.1 | B | | | | | 14.100 | P |
| Lead | | 1.5 | U | | | | | 1.500 | P |
| Magnesium | | 53.8 | B | | | | | 36.700 | P |
| Manganese | | -1.7 | B | | | | | -0.2201 | P |
| Mercury | | | | | | | | 0.100 | AV |
| Nickel | | 0.7 | U | | | | | 0.730 | P |
| Potassium | | 42.1 | U | | | | | 42.100 | P |
| Selenium | | 4.4 | U | | | | | 4.400 | P |
| Silver | | 1.0 | B | | | | | 0.400 | P |
| Sodium | | 45.4 | U | | | | | 45.400 | P |
| Thallium | | 2.6 | U | | | | | 2.600 | P |
| Vanadium | | 0.8 | U | | | | | 0.800 | P |
| Zinc | | 2.4 | B | | | | | 5.815 | P |
| Cyanide | | 0.5 | U | | | | | | AS |

U.S. EPA - CLP
5A
SPIKE SAMPLE RECOVERY

EPA SAMPLE NO.
SKSW51MS1009

Lab Name: GCAL

Lab Code: LA024 Case No.: _____

Contract: _____

Matrix: (soil / water) Water

SAS No.: _____ SDG No.: 204032408

% Solids for Sample: _____

Level (low/med) : _____

Lab Sample ID: 20403240802

Original Lab Sample ID: 20403240803

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| Analyte | Control Limit | | Spiked Sample | | Sample | | Spike Added (SA) | % R | Q | M |
|-----------|---------------|-----|---------------|---|-------------|---|------------------|------|---|----|
| | %R | | Result (SSR) | C | Result (SR) | C | | | | |
| Aluminum | | | 82700 | | 62300 | | 2000 | 1020 | | P |
| Antimony | 75 | 125 | 41.9 | B | 7.3 | B | 100 | 35 | N | P |
| Arsenic | 75 | 125 | 111 | | 50.7 | | 40.0 | 151 | N | P |
| Barium | 75 | 125 | 2580 | | 499 | | 2000 | 104 | | P |
| Beryllium | 75 | 125 | 56.9 | | 4.9 | B | 50.0 | 104 | | P |
| Cadmium | 75 | 125 | 51.6 | | 5.0 | | 50.0 | 93 | | P |
| Chromium | 75 | 125 | 284 | | 72.6 | | 200 | 106 | | P |
| Cobalt | 75 | 125 | 545 | | 59.7 | | 500 | 97 | | P |
| Copper | 75 | 125 | 436 | | 131 | | 250 | 122 | | P |
| Iron | | | 168000 | | 124000 | | 1000 | 4410 | | P |
| Lead | | | 183 | | 122 | | 20.0 | 306 | | P |
| Manganese | | | 8350 | | 5690 | | 500 | 532 | | P |
| Mercury | 75 | 125 | 5.8 | | 0.1 | B | 5.0 | 113 | | AV |
| Nickel | 75 | 125 | 619 | | 116 | | 500 | 101 | | P |
| Selenium | 75 | 125 | 4.4 | U | 4.4 | U | 10.0 | 0 | N | P |
| Silver | 75 | 125 | 46.3 | | 0.4 | U | 50.0 | 93 | | P |
| Thallium | 75 | 125 | 56.7 | | 2.6 | U | 50.0 | 113 | | P |
| Vanadium | 75 | 125 | 614 | | 105 | | 500 | 102 | | P |
| Zinc | 75 | 125 | 1130 | | 490 | | 500 | 128 | N | P |
| Cyanide | 75 | 125 | 87.8 | | 0.8 | B | 100 | 87 | | AS |

Comments:

U.S. EPA - CLP
5A
SPIKE SAMPLE RECOVERY

EPA SAMPLE NO.

SKSW51MS1009 (DISS)

Lab Name: GCAL

Lab Code: LA024 Case No.: _____

Contract: _____

Matrix: (soil / water) Water

SAS No.: _____ SDG No.: 204032408

% Solids for Sample: _____

Level (low/med): _____

Lab Sample ID: 20403240809

Original Lab Sample ID: 20403240810

Concentration Units (ug/L or mg/kg dry weight): ug/L

| Analyte | Control Limit | | Spiked Sample | | Sample Result (SR) | | Spike Added (SA) | % R | Q | M |
|-----------|---------------|-----|---------------|---|--------------------|---|------------------|-----|---|----|
| | 75 | 125 | Result (SSR) | C | Result (SR) | C | | | | |
| Aluminum | 75 | 125 | 2190 | | 25.8 | U | 2000 | 110 | | P |
| Antimony | 75 | 125 | 124 | | 3.7 | U | 100 | 124 | | P |
| Arsenic | 75 | 125 | 51.2 | | 2.9 | U | 40.0 | 128 | N | P |
| Barium | 75 | 125 | 2140 | | 35.8 | B | 2000 | 105 | | P |
| Beryllium | 75 | 125 | 55.8 | | 0.1 | U | 50.0 | 112 | | P |
| Cadmium | 75 | 125 | 52.5 | | 0.2 | U | 50.0 | 105 | | P |
| Chromium | 75 | 125 | 212 | | 1.6 | B | 200 | 105 | | P |
| Cobalt | 75 | 125 | 515 | | 0.4 | B | 500 | 103 | | P |
| Copper | 75 | 125 | 291 | | 4.3 | B | 250 | 115 | | P |
| Iron | 75 | 125 | 1110 | | 14.1 | U | 1000 | 111 | | P |
| Lead | 75 | 125 | 19.0 | | 1.5 | U | 20.0 | 95 | | P |
| Manganese | 75 | 125 | 562 | | 30.0 | | 500 | 106 | | P |
| Mercury | 75 | 125 | 5.8 | | 0.1 | U | 5.0 | 116 | | AV |
| Nickel | 75 | 125 | 519 | | 1.0 | B | 500 | 104 | | P |
| Selenium | 75 | 125 | 4.4 | U | 4.4 | U | 10.0 | 0 | N | P |
| Silver | 75 | 125 | 55.0 | | 0.4 | U | 50.0 | 110 | | P |
| Thallium | 75 | 125 | 50.2 | | 2.6 | U | 50.0 | 100 | | P |
| Vanadium | 75 | 125 | 537 | | 2.2 | B | 500 | 107 | | P |
| Zinc | 75 | 125 | 504 | | 0.6 | U | 500 | 101 | | P |

Comments:

U.S. EPA - CLP
5B
POST DIGEST SPIKE SAMPLE RECOVERY

EPA SAMPLE NO.

SKSW511009PDS

Lab Name: GCAL

Lab Code: LA024 Case No.: _____

Contract: _____

Matrix: (soil / water) Water

SAS No.: _____ SDG No.: 204032408

% Solids for Sample: _____

Level: (low / med) _____

Original Lab Sample ID: 20403240802

Lab Sample ID:: 155382

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| Analyte | Control Limit %R | Spiked Sample Result (SSR) | C | Sample Result (SR) | C | Spike Added (SA) | % R | Q | M |
|-----------|------------------------|----------------------------------|---|-----------------------|---|---------------------|-----|---|---|
| Antimony | | 137 | | 7.3 | B | 120 | 109 | | P |
| Arsenic | | 77.0 | | 50.7 | | 20.0 | 131 | | P |
| Beryllium | | 15.9 | | 4.9 | B | 10.0 | 110 | | P |
| Cadmium | | 14.5 | | 5.0 | | 10.0 | 95 | | P |
| Chromium | | 104 | | 72.6 | | 20.0 | 159 | | P |
| Cobalt | | 156 | | 59.7 | | 100 | 97 | | P |
| Copper | | 186 | | 131 | | 50.0 | 109 | | P |
| Lead | | 128 | | 122 | | 6.0 | 100 | | P |
| Manganese | | 5700 | | 5690 | | 30.0 | 40 | | P |
| Nickel | | 195 | | 116 | | 80.0 | 98 | | P |
| Selenium | | 6.4 | | 4.4 | U | 10.0 | 64 | | P |
| Silver | | 17.6 | | 0.4 | U | 20.0 | 88 | | P |
| Thallium | | 25.8 | | 2.6 | U | 20.0 | 129 | | P |
| Vanadium | | 230 | | 105 | | 100 | 124 | | P |
| Zinc | | 545 | | 490 | | 40.0 | 139 | | P |

Comments:

U.S. EPA - CLP

5B

POST DIGEST SPIKE SAMPLE RECOVERY

EPA SAMPLE NO.

SKSW511009 (DISS)PDS

Lab Name: GCALLab Code: LA024

Case No.: _____

Contract: _____

Matrix: (soil / water) Water

SAS No.: _____

SDG No.: 204032408

% Solids for Sample: _____

Level: (low / med) _____

Original Lab Sample ID: 20403240809

Lab Sample ID:: _____

155384Concentration Units (ug/L or mg/kg dry weight) : ug/L

| Analyte | Control Limit %R | Spiked Sample Result (SSR) | C | Sample Result (SR) | C | Spike Added (SA) | % R | Q | M |
|-----------|------------------------|----------------------------------|---|-----------------------|---|---------------------|-----|---|---|
| Antimony | | 145 | | 3.7 | U | 120 | 121 | | P |
| Arsenic | | 21.2 | | 2.9 | U | 20.0 | 106 | | P |
| Beryllium | | 11.1 | | 0.1 | U | 10.0 | 111 | | P |
| Cadmium | | 10.1 | | 0.2 | U | 10.0 | 101 | | P |
| Chromium | | 22.8 | | 1.6 | B | 20.0 | 106 | | P |
| Cobalt | | 102 | | 0.4 | B | 100 | 101 | | P |
| Copper | | 66.8 | | 4.3 | B | 50.0 | 125 | | P |
| Lead | | 2.8 | B | 1.5 | U | 6.0 | 47 | | P |
| Manganese | | 61.2 | | 30.0 | | 30.0 | 104 | | P |
| Nickel | | 82.4 | | 1.0 | B | 80.0 | 102 | | P |
| Selenium | | 9.4 | | 4.4 | U | 10.0 | 94 | | P |
| Silver | | 22.7 | | 0.4 | U | 20.0 | 113 | | P |
| Thallium | | 13.0 | | 2.6 | U | 20.0 | 65 | | P |
| Vanadium | | 109 | | 2.2 | B | 100 | 106 | | P |
| Zinc | | 40.8 | | 0.6 | U | 40.0 | 102 | | P |

Comments:

U.S. EPA - CLP

6

DUPLICATES

EPA SAMPLE NO

SKSW51DUP1009

Lab Name: GCALLab Code: LA024

SDG No.:

204032408Matrix: (soil / water) Water

% Solids for Sample: _____

Lab Sample ID

20403240805

% Solids for Duplicate: _____

Original Lab Sample

20403240802

Concentration Units (ug/L or mg/kg dry weight) ug/L

| Analyte | Control Limit | Sample (S) | C | Duplicate (D) | C | RPD | Q | M |
|-----------|---------------|------------|---|---------------|---|--------|---|-----|
| Aluminum | 0 - 20 | 62300 | | 88100 | | 34 | * | P ✓ |
| Antimony | - | 7.3 | B | 6.7 | B | 9 | | P |
| Arsenic | 0 - 20 | 50.7 | | 67.0 | | 28 | * | P ✓ |
| Barium | 0 - 200 | 499 | | 714 | | 215.06 | * | P ✓ |
| Beryllium | 0 - 5 | 4.9 | B | 6.7 | | 1.7832 | / | P |
| Cadmium | 0 - 5 | 5.0 | | 7.2 | | 2.1793 | / | P |
| Calcium | 0 - 20 | 427000 | | 603000 | | 34 | * | P ✓ |
| Chromium | 0 - 20 | 72.6 | | 102 | | 34 | * | P ✓ |
| Cobalt | 0 - 50 | 59.7 | | 83.4 | | 23.689 | / | P |
| Copper | 0 - 20 | 131 | | 174 | | 28 | * | P ✓ |
| Iron | 0 - 20 | 124000 | | 173000 | | 33 | * | P ✓ |
| Lead | 0 - 20 | 122 | | 169 | | 32 | * | P ✓ |
| Magnesium | 0 - 20 | 80300 | | 102000 | | 24 | * | P ✓ |
| Manganese | 0 - 20 | 5690 | | 7950 | | 33 | * | P ✓ |
| Mercury | 0 - .2 | 0.1 | B | 0.3 | | .19198 | / | AV |
| Nickel | 0 - 40 | 116 | | 162 | | 45.665 | * | P ✓ |
| Potassium | 0 - 5000 | 12200 | | 17000 | | 4784.0 | / | P |
| Selenium | - | 4.4 | U | 4.4 | U | 0 | | P |
| Silver | - | 0.4 | U | 0.4 | U | 0 | | P |
| Sodium | 0 - 20 | 60200 | | 60100 | | .2 | | P |
| Thallium | - | 2.6 | U | 2.6 | U | 0 | | P |
| Vanadium | 0 - 50 | 105 | | 145 | | 39.768 | | P |
| Zinc | 0 - 20 | 490 | | 679 | | 32 | * | P ✓ |
| Cyanide | - | 0.8 | B | 1.1 | B | 32 | | AS |

U.S. EPA - CLP

6

DUPLICATES

EPA SAMPLE NO

SKSW51DUP1009 (DISS)

Lab Name: GCAL

Lab Code: LA024

SDG No.:

204032408

Matrix: (soil / water) Water

% Solids for Sample:

Lab Sample ID

20403240811

% Solids for Duplicate:

Original Lab Sample

20403240809

Concentration Units (ug/L or mg/kg dry weight) ug/L

| Analyte | Control Limit | Sample (S) | C | Duplicate (D) | C | RPD | Q | M |
|-----------|---------------|------------|---|---------------|---|--------|---|----|
| Aluminum | - | 25.8 | U | 25.8 | U | 0 | | P |
| Antimony | - | 3.7 | U | 3.7 | U | 0 | | P |
| Arsenic | - | 2.9 | U | 2.9 | U | 0 | | P |
| Barium | - | 35.8 | B | 33.0 | B | 8 | | P |
| Beryllium | - | 0.1 | U | 0.1 | U | 0 | | P |
| Cadmium | - | 0.2 | U | 0.2 | U | 0 | | P |
| Calcium | 0 - 20 | 103000 | | 101000 | | 2 | | P |
| Chromium | - | 1.6 | B | 1.4 | B | 13 | | P |
| Cobalt | - | 0.4 | B | 0.4 | B | 0 | | P |
| Copper | - | 4.3 | B | 1.6 | B | 92 | | P |
| Iron | - | 14.1 | U | 14.1 | U | 0 | | P |
| Lead | - | 1.5 | U | 1.5 | U | 0 | | P |
| Magnesium | 0 - 20 | 29700 | | 29100 | | 2 | | P |
| Manganese | 0 - 15 | 30.0 | | 29.7 | | .33702 | | P |
| Mercury | - | 0.1 | U | 0.1 | U | 0 | | AV |
| Nickel | - | 1.0 | B | 1.0 | B | 0 | | P |
| Potassium | - | 2980 | B | 2670 | B | 11 | | P |
| Selenium | - | 4.4 | U | 4.4 | U | 0 | | P |
| Silver | - | 0.4 | U | 0.4 | U | 0 | | P |
| Sodium | 0 - 20 | 59200 | | 57300 | | 3 | | P |
| Thallium | - | 2.6 | U | 2.6 | U | 0 | | P |
| Vanadium | - | 2.2 | B | 1.7 | B | 26 | | P |
| Zinc | - | 0.6 | U | 0.6 | U | 0 | | P |

U.S. EPA - CLP
7
LABORATORY CONTROL SAMPLE

Lab Name: GCAL Contract: _____
 Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204032408
 Solid LCS Source: _____ Lab Sample ID: 154787
 Aqueous LCS Source: 310006 HIGH PURITY~323003 HIGH

| Analyte | Aqueous (ug/L) | | | Solid (mg/kg) | | | |
|-----------|----------------|-------|-----|---------------|-------|---|-----|
| | True | Found | % R | True | Found | C | % R |
| Aluminum | 2000 | 2150 | 107 | | | | |
| Antimony | 500 | 572 | 114 | | | | |
| Arsenic | 2000 | 2190 | 109 | | | | |
| Barium | 2000 | 2070 | 104 | | | | |
| Beryllium | 50.0 | 54.7 | 109 | | | | |
| Cadmium | 50.0 | 51.1 | 102 | | | | |
| Calcium | 12500 | 12900 | 103 | | | | |
| Chromium | 200 | 205 | 103 | | | | |
| Cobalt | 500 | 504 | 101 | | | | |
| Copper | 250 | 280 | 112 | | | | |
| Iron | 1000 | 1110 | 111 | | | | |
| Lead | 500 | 523 | 105 | | | | |
| Magnesium | 12500 | 12500 | 100 | | | | |
| Manganese | 500 | 523 | 105 | | | | |
| Nickel | 500 | 512 | 102 | | | | |
| Potassium | 12500 | 12400 | 99 | | | | |
| Selenium | 2000 | 2200 | 110 | | | | |
| Silver | 50.0 | 54.0 | 108 | | | | |
| Sodium | 12500 | 13200 | 106 | | | | |
| Thallium | 2000 | 2100 | 105 | | | | |
| Vanadium | 500 | 526 | 105 | | | | |
| Zinc | 500 | 502 | 100 | | | | |

U.S. EPA - CLP
7
LABORATORY CONTROL SAMPLE

Lab Name: GCAL Contract: _____
 Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204032408
 Solic LCS Source: _____ Lab Sample ID: 154791
 Aqueous LCS Source: 310006 HIGH PURITY~323003 HIGH

| Analyte | Aqueous (ug/L) | | | Solid (mg/kg) | | | |
|-----------|----------------|-------|-----|---------------|-------|---|-----|
| | True | Found | % R | True | Found | C | % R |
| Aluminum | 2000 | 2130 | 106 | | | | |
| Antimony | 500 | 578 | 116 | | | | |
| Arsenic | 2000 | 2190 | 109 | | | | |
| Barium | 2000 | 2070 | 103 | | | | |
| Beryllium | 50.0 | 54.6 | 109 | | | | |
| Cadmium | 50.0 | 51.2 | 102 | | | | |
| Calcium | 12500 | 13000 | 104 | | | | |
| Chromium | 200 | 205 | 102 | | | | |
| Cobalt | 500 | 503 | 101 | | | | |
| Copper | 250 | 279 | 112 | | | | |
| Iron | 1000 | 1050 | 105 | | | | |
| Lead | 500 | 520 | 104 | | | | |
| Magnesium | 12500 | 12600 | 101 | | | | |
| Manganese | 500 | 520 | 104 | | | | |
| Nickel | 500 | 510 | 102 | | | | |
| Potassium | 12500 | 12500 | 100 | | | | |
| Selenium | 2000 | 2190 | 110 | | | | |
| Silver | 50.0 | 53.8 | 108 | | | | |
| Sodium | 12500 | 13400 | 107 | | | | |
| Thallium | 2000 | 2090 | 105 | | | | |
| Vanadium | 500 | 527 | 105 | | | | |
| Zinc | 500 | 495 | 99 | | | | |

U.S. EPA - CLP

9

ICP SERIAL DILUTIONS

EPA SAMPLE NO.

SKSW511009SD

Lab Name: GCALLab Code: LA024

Case No. _____

SDG No.: 204032408Matrix: (soil / water) WaterLab Sample ID 155381Original Lab Sample ID 20403240802Concentration Units: ug/L

| Analyte | Initial Sample Result (I) | C | Serial Dilution Result (S) | C | % Difference | Q | M |
|-----------|---------------------------------|---|----------------------------------|---|--------------|---|-----|
| Aluminum | 62300 | | 131000 | | 110.3 | E | P ✓ |
| Antimony | 7.3 | B | 18.5 | U | 153.4 | | P |
| Arsenic | 50.7 | | 69.6 | | 37.3 | | P |
| Barium | 499 | | 837 | B | 67.7 | E | P ✓ |
| Beryllium | 4.9 | B | 7.4 | B | 51.0 | | P |
| Cadmium | 5.0 | | 4.4 | B | 12.0 | | P |
| Calcium | 427000 | | 462000 | | 8.2 | | P |
| Chromium | 72.6 | | 135 | | 86.0 | E | P ✓ |
| Cobalt | 59.7 | | 61.0 | B | 2.2 | | P |
| Copper | 131 | | 143 | | 9.2 | | P |
| Iron | 124000 | | 148000 | | 19.4 | E | P ✓ |
| Lead | 122 | | 121 | | 0.8 | | P |
| Magnesium | 80300 | | 90500 | | 12.7 | E | P ✓ |
| Manganese | 5690 | | 6180 | | 8.6 | | P |
| Nickel | 116 | | 132 | B | 13.8 | E | P ✓ |
| Potassium | 12200 | | 41600 | | 241.0 | E | P ✓ |
| Selenium | 4.4 | U | 22.0 | U | | | P |
| Silver | 0.4 | U | 2.0 | U | | | P |
| Sodium | 60200 | | 61500 | | 2.2 | | P |
| Thallium | 2.6 | U | 13.0 | U | | | P |
| Vanadium | 105 | | 221 | B | 110.5 | E | P ✓ |
| Zinc | 490 | | 489 | | 0.2 | | P |

U.S. EPA - CLP
9
ICP SERIAL DILUTIONS

EPA SAMPLE NO.

SKSW511009 (DISS)SD

Lab Name: GCAL

Lab Code: LA024

Case No.

SDG No.: 204032408

Matrix: (soil / water) Water

Lab Sample ID 155383

Original Lab Sample ID 20403240809

Concentration Units: ug/L

| Analyte | Initial Sample Result (I) | C | Serial Dilution Result (S) | C | % Difference | Q | M |
|-----------|---------------------------------|---|----------------------------------|---|--------------|---|---|
| Aluminum | 25.8 | U | 129 | U | | | P |
| Antimony | 3.7 | U | 18.5 | U | | | P |
| Arsenic | 2.9 | U | 14.5 | U | | | P |
| Barium | 35.8 | B | 34.8 | B | 2.8 | | P |
| Beryllium | 0.1 | U | 0.5 | U | | | P |
| Cadmium | 0.2 | U | 1.0 | U | | | P |
| Calcium | 103000 | | 104000 | | 1.0 | | P |
| Chromium | 1.6 | B | 4.0 | U | 150.0 | | P |
| Cobalt | 0.4 | B | 2.0 | U | | | P |
| Copper | 4.3 | B | 6.0 | U | 39.5 | | P |
| Iron | 14.1 | U | 70.5 | U | | | P |
| Lead | 1.5 | U | 7.5 | U | | | P |
| Magnesium | 29700 | | 29700 | | 0.0 | | P |
| Manganese | 30.0 | | 23.4 | B | 22.0 | E | P |
| Nickel | 1.0 | B | 4.3 | B | 330.0 | | P |
| Potassium | 2980 | B | 2790 | B | 6.4 | | P |
| Selenium | 4.4 | U | 22.0 | U | | | P |
| Silver | 0.4 | U | 2.0 | U | | | P |
| Sodium | 59200 | | 57100 | | 3.5 | | P |
| Thallium | 2.6 | U | 13.0 | U | | | P |
| Vanadium | 2.2 | B | 4.0 | U | 81.8 | | P |
| Zinc | 0.6 | U | 3.0 | U | | | P |

U.S. EPA - CLP

10

INSTRUMENT DETECTION LIMITS

Lab Name: GCALLab Code: LA024

Case No.: _____

SAS No.: _____

SDG No.: 204032408Instrument ID: ICP5Study Date: 04/30/03

| <i>Analyte</i> | <i>Wavelength (nm)</i> | <i>Background</i> | <i>CRDL (ug/L)</i> | <i>IDL (ug/L)</i> | <i>M</i> |
|----------------|----------------------------|-------------------|------------------------|-----------------------|----------|
| Aluminum | 308.210 | | 200 | 25.8 | P |
| Antimony | 206.830 | | 60 | 3.7 | P |
| Arsenic | 193.700 | | 10 | 2.9 | P |
| Barium | 233.520 | | 200 | .3 | P |
| Beryllium | 313.100 | | 5 | .1 | P |
| Cadmium | 214.430 | | 5 | .2 | P |
| Calcium | 315.880 | | 5000 | 7.5 | P |
| Chromium | 267.710 | | 10 | .8 | P |
| Cobalt | 228.610 | | 50 | .4 | P |
| Copper | 324.750 | | 25 | 1.2 | P |
| Iron | 259.940 | | 100 | 14.1 | P |
| Lead | 220.350 | | 3 | 1.5 | P |
| Magnesium | 279.080 | | 5000 | 36.7 | P |
| Manganese | 257.610 | | 15 | .2 | P |
| Nickel | 231.600 | | 40 | .7 | P |
| Potassium | 766.480 | | 5000 | 42.1 | P |
| Selenium | 196.030 | | 5 | 4.4 | P |
| Silver | 328.060 | | 10 | .4 | P |
| Sodium | 589.580 | | 5000 | 45.4 | P |
| Thallium | 190.800 | | 10 | 2.6 | P |
| Vanadium | 290.880 | | 50 | .8 | P |
| Zinc | 213.860 | | 20 | .6 | P |

U.S. EPA - CLP
14
ANALYSIS RUN LOG

Lab Name: GCAL
Lab Code: LA024 Case No.: _____
Instrument ID Number: ICP5
Start Date: 03/29/04

Contract: _____
SAS No.: _____ SDG No.: 204032408
Method Type: P
End Date: 03/29/04

Analyte Symbols

| EPA Sample No. | D/F | Time | % R | Al | Sb | As | Ba | Be | Cd | Ca | Cr | Co | Cu | Fe | Pb | Mg | Mn | Hg | Ni | K | Se | Ag | Na | Tl | V | Zn | Cn |
|----------------------|-----|------|-----|----|----|----|----|----|----|----|----|----|----|----|----|----|----|----|----|---|----|----|----|----|---|----|----|
| ICV | 1 | 1155 | | | X | X | | X | X | X | X | X | X | X | X | X | X | | X | | X | X | | X | X | X | |
| ICV2 | 1 | 1201 | | X | | | X | | | | | | | | | | | | | X | | | X | | | | |
| ICB | 1 | 1208 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| CRDL | 1 | 1215 | | | X | X | | X | X | | X | X | X | | X | | X | | X | | X | X | | X | X | X | |
| ICSA | 1 | 1222 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| ICSAB | 1 | 1228 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| CCV | 1 | 1234 | | X | | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| CCV2 | 1 | 1240 | | | X | | | | | | | | | | | | | | | | | | | | | | |
| CCB | 1 | 1248 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| MB154786 | 1 | 1254 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKSW511009 | 1 | 1301 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKSW51DUP1009 | 1 | 1308 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKSW511009SD | 5 | 1314 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKSWEB1009 | 1 | 1320 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKSW521009 | 1 | 1327 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKSW531009 | 1 | 1333 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKSW51MS1009 | 1 | 1340 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKSW511009PDS | 1 | 1346 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| LCS154787 | 1 | 1352 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| CCV | 1 | 1358 | | X | | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| CCV2 | 1 | 1405 | | | X | | | | | | | | | | | | | | | | | | | | | | |
| CCB | 1 | 1412 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| MB154790 | 1 | 1419 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKSW511009 (DISS) | 1 | 1425 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKSW51DUP1009 (DISS) | 1 | 1432 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKSW511009 (DISS)SD | 5 | 1439 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKSWEB1009 (DISS) | 1 | 1446 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |

U.S. EPA - CLP
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ANALYSIS RUN LOG

Lab Name: GCAL
Lab Code: LA024 Case No.: _____
Instrument ID Number: ICP5
Start Date: 03/29/04

Contract: _____
SAS No.: _____ SDG No.: 204032408
Method Type: P
End Date: 03/29/04

Analyte Symbols

| EPA Sample No. | D/F | Time | % R | Al | Sb | As | Ba | Be | Cd | Ca | Cr | Co | Cu | Fe | Pb | Mg | Mn | Hg | Ni | K | Se | Ag | Na | Tl | V | Zn | Cn |
|----------------------|-----|------|-----|----|----|----|----|----|----|----|----|----|----|----|----|----|----|----|----|---|----|----|----|----|---|----|----|
| SKSW521009 (DISS) | 1 | 1453 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKSW531009 (DISS) | 1 | 1500 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| CRDL | 1 | 1506 | | | X | X | | X | X | | X | X | X | | X | | X | | X | | X | X | | X | X | X | |
| ICSA | 1 | 1513 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| ICSAB | 1 | 1519 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| CCV | 1 | 1525 | | X | | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| CCV2 | 1 | 1532 | | | X | | | | | | | | | | | | | | | | | | | | | | |
| CCB | 1 | 1539 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKSW51MS1009 (DISS) | 1 | 1546 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKSW511009 (DISS)PDS | 1 | 1552 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| LCS154791 | 1 | 1559 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| CRDL | 1 | 1605 | | | X | X | | X | X | | X | X | X | | X | | X | | X | | X | X | | X | X | X | |
| ICSA | 1 | 1612 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| ICSAB | 1 | 1618 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| CCV | 1 | 1624 | | X | | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| CCV2 | 1 | 1630 | | | X | | | | | | | | | | | | | | | | | | | | | | |
| CCB | 1 | 1637 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |

DATA VALIDATION REPORT
FOR
SKINNER LANDFILL SITE
EARTH TECH: PROJECT NUMBER 54280
LABORATORY REPORT NUMBER 204030804
PROJECT MANAGER: Ron Rolker
Date: May 6, 2004
Revised Report Dated: September 23, 2005
Data Validator: Mark Kromis

APPENDIX C LIST OF ACRONYMS

| | |
|--------|--|
| BFB | Bromofluorobenzene |
| CC | Continuing Calibration |
| CCV | Continuing Calibration Verification |
| CCB | Continuing Calibration Blanks |
| CLP | Contract Laboratory Program |
| CRDL | Contract Required Detection Limit |
| DFTPP | Decafluorotriphenylphosphine |
| GC/MS | Gas Chromatograph/Mass Spectrometer |
| IC | Initial Calibration |
| ICB | Initial Calibration Blank |
| IDL | Instrument Detection Limit |
| ICP | Inductively Coupled Plasma |
| ICS | Interference Check Sample |
| ICV | Initial Calibration Verification |
| ILM | Inorganic Analysis Multi-Media Multi-Concentration |
| INDAM | Individual A Mixture |
| INDBM | Individual B Mixture |
| mg/L | milligrams per liter |
| MS/MSD | Matrix Spike/Matrix Spike Duplicate |
| OLC | Organic Analysis Low Concentration |
| OLM | Organic Analysis Multi-Media Multi-Concentration |
| %D | Percent Difference |
| % RSD | Percent Relative Standard Deviation |
| PB | Preparation Blanks |
| QC | Quality Control |
| RF | Response Factor |
| RPD | Relative Percent Difference |
| RRF | Relative Response Factor |
| SDG | Sample Delivery Group |
| SOW | Statement of Work |
| µg/L | micrograms per liter |
| US EPA | United States Environmental Protection Agency |
| VOC | Volatile Organic Compounds |
| VTSR | Validated Time of Sample Receipt |

DATA VALIDATION SUMMARY – SAMPLE DELIVERY GROUP 204030804 INORGANICS

Validation of the inorganics data, as prepared by Gulf Coast Analytical Laboratories (GCAL) for the samples collected from the Skinner Landfill site in March 2004, was conducted by Earth Tech using the National Functional Guidelines for Inorganic Data Review, (US EPA, February, 1994), as appropriate. The results were reported by GCAL under Sample Delivery Group (SDG) 204030804.

| GCAL # | Sample Description |
|--------------|-----------------------|
| 204030804-01 | SKSWD031009 |
| 204030804-02 | SKSWD03D1009 |
| 204030804-03 | SKSWD03MS1009 |
| 204030804-05 | SKSWD03DUP1009 |
| 204030804-06 | SKSWDEB1009 |
| 204030804-08 | SKSWD031009 (DISS) |
| 204030804-09 | SKSWD03D1009 (DISS) |
| 204030804-10 | SKSWD03MS1009 (DISS) |
| 204030804-11 | SKSWD03DUP1009 (DISS) |
| 204030804-12 | SKSWDEB1009 (DISS) |

INTRODUCTION

Analyses of metals were performed according to Contract Laboratory Program (CLP)-Inorganic Analysis Multi-media Multi-concentration ILM04.1 Statement of Work (SOW). Results of the sample analyses are reported by the laboratory as either qualified or unqualified. Unqualified results mean that the reported values maybe used without reservation. Various qualifier codes are used by the laboratory to denote specific information regarding the analytical results.

The data validation process is intended to evaluate the data on a technical basis. The data package also was subjected to an internal laboratory quality review prior to submission to Earth Tech for data validation.

During the validation process, laboratory-qualified and unqualified data are verified against all available supporting documentation. Based on this evaluation, qualifier codes may be added, deleted or modified by the data user. Final results are therefore, either qualified or unqualified. Validator-qualified results are annotated with the following codes in accordance with the Functional Guidelines:

- U The constituent was analyzed for, but was not detected above the level of the associated analytical reporting limit. The associated value is either the sample quantitation limit or the sample detection limit.
- J The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample.

- UJ The analyte was not detected above the reported sample quantitation limit. However, the reported quantitation limit is approximate and may or may not represent the actual limit of quantitation necessary to accurately and precisely measure the analyte in the sample.
- R The sample results are rejected due to serious deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte cannot be verified.

Details of the inorganics data validation findings and conclusions are provided in the following sections of this report:

1. Holding Times
2. Calibration
 - A. Initial Calibration (IC)
 - B. Continuing Calibration (CC)
3. Blanks
4. Inductively Coupled Plasma (ICP) Interference Check Sample
5. Laboratory Control Sample (LCS)
6. Duplicate Analysis
7. Spike Sample Analysis
8. ICP Serial Dilution
9. System Performance
10. Documentation
11. Overall Assessment

1. HOLDING TIMES

All samples for inorganics analyses were analyzed within the 180-day holding time for preserved aqueous samples. Mercury analyses were conducted within the 28-day holding time for aqueous samples undergoing CLP protocol. Cyanide analyses were conducted within the 14-day holding time. The cooler temperature upon receipt at the laboratory was within the recommended temperature of 4°C +/- 2°C.

2. CALIBRATION

A. Initial Calibration

The percent recoveries for the Initial Calibration Verification (ICV) standard were within Quality Control (QC) limits for all constituents.

B. Continuing Calibration

The percent recoveries for the Continuing Calibration Verification (CCV) standard were within QC limits for all constituents.

3. BLANKS

The Initial Calibration Blank (ICB), Continuing Calibration Blanks (CCB) and Preparation Blanks (PB) were analyzed at the appropriate frequencies. No constituents were detected in the ICB, CCB, PB, Equipment blanks above the corresponding Contract Required Detection Limit (CRDL).

4. ICP INTERFERENCE CHECK SAMPLE

Results for the ICP analysis of the Interference Check Sample (ICS) solution AB were within 20% of the true value.

5. LABORATORY CONTROL SAMPLES

Recoveries were within the control limit (80-120%) for all constituents.

6. DUPLICATE ANALYSIS

The Relative Percent Difference (RPD) between the sample and duplicate results were within the acceptance criteria for all target compounds.

7. SPIKE SAMPLE ANALYSIS

The laboratory used sample SKSWD031009 for the matrix spike sample. The MS percent recoveries were within the acceptance criteria (75%-125%) with the exception of Selenium in the total (0%) and dissolved (50%) fractions. As per the National Functional Guidelines: if the percent recovery is greater than 30% and less than 74% qualify detected results for that analyte with "J" and non-detected results with "UJ". If the percent recovery is less than 30% qualify detected results for that analyte with "J" and non-detected results with "R".

8. ICP SERIAL DILUTION

As noted in the National Functional Guidelines: If the analyte concentration is at least 50 times above the IDL, its serial dilution analysis must then agree within 10% of the original determination after corrected for dilution. The serial dilution is performed to determine whether any significant chemical or physical interference's exist due to matrix effects. The serial dilution percent differences were within the acceptance criteria for all target analytes with the exception of Aluminum and Zinc in the total fraction. As per the National Functional Guidelines, if the required 10% difference criterion is not met then qualify the associated results as estimated with "J".

9. SYSTEM PERFORMANCE

The analytical system appears to have been working well at the time of these analyses, based on the evaluation of the raw data.

10. DOCUMENTATION

The documentation appeared accurate and in order with the exception of Manganese associated with the dissolved fraction. GCAL qualified the Manganese results with an "E" to indicate that the serial dilution percent difference criteria had been exceeded. The initial concentration of Manganese was not greater than 50 times the IDL therefore the analyte did not need to be flagged. The data validator crossed the "E" qualifier out with a single line and initialed and dated the correction.

11. OVERALL ASSESSMENT

The percent recoveries for Lead in the Contract Required Detection Limit (CRDL) standards were 101%, 78%, and 123%.

The percent recoveries for Copper in the Contract Required Detection Limit (CRDL) standards were 63.0, 62.9%, and 69.1%.

The percent recoveries for Selenium in the Contract Required Detection Limit (CRDL) standards were 62%, 64%, and 84%.

The percent recoveries for Zinc in the Contract Required Detection Limit (CRDL) standards were 76%, 75%, and 81%.

If the CRDL is greater than 120% then detected results greater than the IDL but less than two times the CRDL are qualified as estimated with "J". If the CRDL is below 80% then detected results are qualified as estimated with "J" and the non-detected results were qualified with "UJ".

The results are acceptable with the validator-added qualifiers.

**DATA VALIDATION SUMMARY – SAMPLE DELIVERY GROUP 204030804
SEMIVOLATILE ORGANICS**

Validation of the Gas Chromatograph/Mass Spectrometer (GC/MS) semi-volatile organics data, as prepared by Gulf Coast Analytical Laboratories (GCAL) for the samples collected from the Skinner Landfill site in March 2004, was conducted by Earth Tech using the National Functional Guidelines for Organic Data Review, (US EPA, October, 1999) as appropriate. The results were reported by GCAL under SDG 204030804.

| GCAL # | Sample Description |
|--------------|--------------------|
| 204030804-01 | SKSWD031009 |
| 204030804-02 | SKSWD03D1009 |
| 204030804-03 | SKSWD03MS1009 |
| 204030804-04 | SKSWD03MSD1009 |
| 204030804-06 | SKSWDEB1009 |

INTRODUCTION

Analyses were performed according to CLP-Organic Analysis Multi-Media, Multi-Concentration OLM04.2 SOW. Results of the sample analyses are reported by the laboratory as either qualified or unqualified. Unqualified results mean that the reported values may be used without reservation. Various data qualifier codes are used by the laboratory to denote specific information regarding the analytical results. The data validation process is intended to evaluate the data on a technical basis. The data package also was subjected to an internal laboratory quality review prior to submission to Earth Tech for data validation.

During the validation process, laboratory-qualified and unqualified data are verified against all available supporting documentation. Based on this evaluation, qualifier codes may be added, deleted or modified by the data user. Final results are therefore, either qualified or unqualified. Validator-qualified results are annotated with the following codes in accordance with the Functional Guidelines:

- U The constituent was analyzed for, but was not detected above the level of the associated analytical reporting limit. The associated value is either the sample quantitation limit or the sample detection limit.
- J The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample.
- UJ The analyte was not detected above the reported sample quantitation limit. However, the reported quantitation limit is approximate and may or may not represent the actual limit of quantitation necessary to accurately and precisely measure the analyte in the sample.

- R The sample results are rejected due to serious deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte cannot be verified.

Details of the semivolatile data validation findings and conclusions are provided in the following sections of this report:

1. Holding Times
2. GC/MS Tuning
3. Calibration
 - A. IC
 - B. CC
4. Blanks
5. System Monitoring Compound Recovery
6. MS/MSD
7. Internal Standards Performance
8. Compound Identification
9. Constituent Quantitation and Reported Detection Limits
10. System Performance
11. Documentation
12. Overall Assessment

1. HOLDING TIMES

All samples were initially extracted within the seven-day technical holding time and the five-day VTSR method holding time. The cooler temperature upon receipt at the laboratory was within the recommended temperature of 4°C +/- 2°C.

2. GC/MS TUNING

The samples were analyzed on a single GC/MS system, identified as MSSV2. One decafluorotriphenylphosphine (DFTPP) tune was run representing the shift in which the standards and samples were analyzed. The DFTPP tune is acceptable.

3. CALIBRATION

A. Initial Calibration

One IC dated 3/19/04 was analyzed in support of the semivolatile sample analyses. Documentation of the IC was present in the data package, and the Relative Response Factor (RRF), as well as percent % RSD values were accurately reported for all target compounds. The criteria employed for technical data review purposes are different than those used in the method. The laboratory must meet a minimum RRF of 0.01; however, for data review purposes, a RRF criterion of "greater than or equal to 0.05" is applied to all semi-volatile compounds. The RRF's and the average RRF were within the acceptance criteria specified in the method for all reported analytes. The %RSD's were within the acceptance criteria specified in the method for all target analytes with the exception of Di-n-butylphthalate (32.0%), Di-n-octylphthalate (30.3%) and Diethylphthalate (41.8%). The lowest point of the calibration curve was dropped for Di-n-butylphthalate, Di-n-octylphthalate, and Diethylphthalate and the %RSD were recalculated. The recalculated %RSD were within the acceptance criteria of less than 30%. Di-n-butylphthalate, Di-n-octylphthalate, and Diethylphthalate were not detected in the associated samples therefore data qualification was not required.

B. Continuing Calibration

One CC dated 3/19/04 was analyzed in support of the semivolatile sample analyses reported in the data submissions. The RRF's for the CC dated 3/19/04 were within the acceptance criteria. The percent difference (%D) between the average RRF's and the CC Response Factors for the CC dated 3/19/04 were within the acceptance criteria with the exception the %D for Naphthalene. As per the National Functional Guidelines, if the %D exceeds the acceptance criteria qualify detected results for that analyte with "J" and non-detected results for that analyte with "UJ".

4. BLANKS

One laboratory semivolatile method blank and equipment blank were analyzed with this SDG. The results are summarized below.

Method Blank (0308SBLK)

There were not target analytes detected in method blank 0308SBLK.

Equipment Blank (SKSWDEB1009)

There were not target analytes detected in method blank SKSWDEB1009 collected on 3/2/04.

5. SYSTEM MONITORING COMPOUND RECOVERY

All reported semivolatile system monitoring compounds were recovered within acceptable control limits.

6. MATRIX SPIKE/MATRIX SPIKE DUPLICATE (MS/MSD)

Sample SKSWD031009 was submitted for MS/MSD analysis. The MS/MSD percent recoveries were within the acceptance criteria with the exception of the 4-Nitorphenol. The %RPD between the MS/MSD were within the acceptance criteria. As per the National Functional Guidelines, no action is taken on MS/MSD data alone.

7. INTERNAL STANDARDS PERFORMANCE

Internal standard areas and retention times were within acceptable limits for the reported semivolatile sample analyses.

8. COMPOUND IDENTIFICATION

All reported semivolatile constituents were correctly identified with supporting chromatograms present in the data package.

9. CONSTITUENT QUANTITATION AND REPORTED DETECTION LIMITS

Constituent quantitations were correctly calculated and reported for semivolatile constituents.

10. SYSTEM PERFORMANCE

The analytical system appears to have been working well at the time of these analyses, based on the evaluation of the raw data submitted for review.

11. DOCUMENTATION

The documentation appeared accurate and in order with the exception of the following: GCAL inadvertently left the Date Extracted off of the CLP Form 1's therefore the data validator inserted the Date Extracted on the CLP Form 1's.

12. OVERALL ASSESSMENT

The results are acceptable with the validator-added qualifiers.

**DATA VALIDATION SUMMARY – SAMPLE DELIVERY GROUP 204030804
VOLATILE ORGANIC**

Validation of the GC/MS volatile organics data, as prepared by Gulf Coast Analytical Laboratories (GCAL) for the samples collected from the Skinner Landfill site in March 2004, was conducted by Earth Tech using the National Functional Guidelines for Organic Data Review, (US EPA, October, 1999), as appropriate. The results were reported by GCAL under SDG 204030804.

| GCAL # | Sample Description |
|--------------|--------------------|
| 204030804-01 | SKSWD031009 |
| 204030804-02 | SKSWD03D1009 |
| 204030804-03 | SKSWD03MS1009 |
| 204030804-04 | SKSWD03MSD1009 |
| 204030804-06 | SKSWDEB1009 |
| 204030804-07 | Trip Blank |

INTRODUCTION

Analyses were performed according to CLP-Organic Analysis Low Concentration OLC02.0 SOW. Results of the sample analyses are reported by the laboratory as either qualified or unqualified. Unqualified results mean that the reported values may be used without reservation. Various qualifier codes are used by the laboratory to denote specific information regarding the analytical results. The data validation process is intended to evaluate the data on a technical basis. The data package also was subjected to an internal laboratory quality review prior to submission to Earth Tech for data validation.

During the validation process, laboratory-qualified and unqualified data are verified against all available supporting documentation. Based on this evaluation, qualifier codes may be added, deleted or modified by the data user. Final results are therefore, either qualified or unqualified. Validator-qualified results are annotated with the following codes in accordance with the Functional Guidelines:

- U The constituent was analyzed for, but was not detected above the level of the associated analytical reporting limit. The associated value is either the sample quantitation limit or the sample detection limit.
- J The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample.
- UJ The analyte was not detected above the reported sample quantitation limit.

However, the reported quantitation limit is approximate and may or may not represent the actual limit of quantitation necessary to accurately and precisely measure the analyte in the sample.

- R The sample results are rejected due to serious deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte cannot be verified.

The volatiles data validation findings and conclusions are provided in the following sections of this report:

1. Holding Times
2. GC/MS Tuning
3. Calibration
 - A. IC
 - B. CC
4. Blanks
5. System Monitoring Compound Recovery
6. MS/MSD
7. Laboratory Control Sample
8. Internal Standards Performance
9. Compound Identification
10. Constituent Quantitation and Reported Detection Limits
11. System Performance
12. Documentation
13. Overall Assessment

1. **HOLDING TIMES**

All samples for Volatile Organic Compounds (VOC) analyses were analyzed within the 14-day technical holding time and the 10-day VTSR method holding time. The cooler temperature upon receipt at the laboratory was within the recommended temperature of 4°C +/- 2°C.

2. GC/MS TUNING

All samples were analyzed on a single GC/MS system, identified as MSV2. One bromofluorobenzene (BFB) tune was run. The BFB tune is acceptable.

3. CALIBRATION

A. Initial Calibration

One IC dated 3/13/04 was analyzed on Instrument MSV2 in support of the volatile sample analyses reported in the data submissions. Documentation of the IC standards was present in the data package, and RRF's as well as %RSD values were accurately reported. The criteria employed for technical data review purposes are different than those used in the method. The laboratory must meet a minimum RRF of 0.01; however, for data review purposes, a RRF criterion of "greater than or equal to 0.05" is applied to all volatile compounds. The %RSD's were within the acceptance criteria specified in the method for all target analytes. The RRF's and the average RRF for the IC dated 3/13/04 were within the acceptance criteria specified in the method for all target analytes with the exception of Acetone. As per the National Functional Guidelines, if any initial calibration RRF is less than 0.05, qualify positive results that have acceptable mass spectral identification with "J", using professional judgement, and non-detected analytes as unusable (R).

B. Continuing Calibration

One CC dated 3/13/04 was analyzed on instrument MSV2 in support of the volatile sample analyses reported in the data submissions. The percent difference (%D) between the average RRF's and the CC RF's were within the acceptance criteria for all target analytes. The CC RRF's were within the acceptance criteria specified in the method for all target analytes with the exception of Acetone. The Acetone results were previously qualified under section 3A above.

4. BLANKS

One laboratory volatile method blank, storage blank, Trip Blank, and an Equipment Blank were analyzed with this SDG. The results are summarized below.

V2BLK01 (MB151783)

Methylene chloride, Chloroform and 1,3-Dichlorobenzene were detected at concentrations of 0.12 ppb, 0.38 ppb, and 0.034 ppb respectively in the method blank analyzed on 3/13/04.

Storage Blank (VHBLK01)

Methylene chloride (0.15 ppb) and 1,4-Dichlorobenzene (0.21 ppb) were detected in the storage blank analyzed on 3/13/04.

Trip Blank

Methylene chloride was detected at a concentration of 0.47 ppb in the Trip Blank submitted for the sampling event that occurred on 3/2/04. The Methylene chloride detected in the trip blank was mitigated by the presence of Methylene chloride in the associated method blank.

Equipment Blank (SKSWDEB1009)

Ethylbenzene (0.022 ppb), Methylene chloride (0.4 ppb), Styrene (0.06 ppb) Toluene (0.43 ppb) and total Xylenes (0.11 ppb) were detected in the Equipment Blank collected on 3/2/04. The analytes Ethylbenzene, Styrene, Toluene, and total Xylenes were not detected in the associated samples therefore no data qualification was not required. The Methylene chloride detected in the equipment blank was mitigated by the presence of Methylene chloride in the associated method blank.

5. SYSTEM MONITORING COMPOUND RECOVERY

All reported volatile system monitoring compounds were recovered within acceptable control limits for all samples.

6. MATRIX SPIKE/MATRIX SPIKE DUPLICATE

Sample SKSWD031009 was submitted for MS/MSD analysis. The MS/MSD percent recoveries and %RPD between the MS/MSD were within the acceptance criteria.

7. LABORATORY CONTROL SAMPLE

A LCS was analyzed in conjunction with this SDG. Recoveries were within the control limit for all constituents.

8. INTERNAL STANDARDS PERFORMANCE

Internal Standard areas and retention times were within acceptable limits for the reported volatile sample analyses.

9. COMPOUND IDENTIFICATION

All reported VOCs were correctly identified with supporting chromatograms present in the data package.

10. CONSTITUENT QUANTITATION AND REPORTED DETECTION LIMITS

Constituent quantitations were correctly calculated and reported for VOCs with the exception of Ethylbenzene. The Ethylbenzene standard and detected results were originally quantitated using the incorrect quantitation ion (GCAL used 106 instead of 91).

GCAL corrected the mistake and re-submitted the corrected pages that were affected in the laboratory report. The overall effect had no impact in the final result for Ethylbenzene.

11. SYSTEM PERFORMANCE

The analytical system appears to have been working well at the time of these analyses, based on the evaluation of the raw data.

12. DOCUMENTATION

The documentation appeared accurate and in order with the exception of the following: GCAL inadvertently left the "B" qualifier off of the CLP Form 1's for Methylene chloride therefore the data validator inserted a "B" qualifier in the "Q" column of the CLP Form 1's. The "B" qualifier indicates that the analyte was detected in the associated method blank.

13. OVERALL ASSESSMENT

The Acetone detected in sample SKSWD03D1009 could be do to low level contamination because Acetone is a common laboratory contaminant and the fact that Acetone was not detected in the associated duplicate sample. The results are acceptable with the validator-added qualifiers.

DATA VALIDATION SUMMARY - SAMPLE DELIVERY GROUP 204030804 PESTICIDES

Validation of the Gas Chromatography (GC) pesticides data, as prepared by Gulf Coast Analytical Laboratories (GCAL) for the samples collected from the Skinner Landfill site in March 2004, was conducted by Earth Tech using the National Functional Guidelines for Organic Data Review, (US EPA, October, 1999), as appropriate. The results were reported by GCAL under SDG 204030804.

| GCAL # | Sample Description |
|--------------|--------------------|
| 204030804-01 | SKSWD031009 |
| 204030804-02 | SKSWD03D1009 |
| 204030804-03 | SKSWD03MS1009 |
| 204030804-04 | SKSWD03MSD1009 |
| 204030804-06 | SKSWDEB1009 |

INTRODUCTION

Analyses were performed according to CLP-Organic Analysis Multi-Media, Multi-Concentration OLM04.2 SOW. Results of the sample analyses are reported by the laboratory as either qualified or unqualified. Unqualified results mean that the reported values may be used without reservation. Various qualifier codes are used by the laboratory to denote specific information regarding the analytical results.

The data validation process is intended to evaluate the data on a technical basis. The data package also was subjected to an internal laboratory quality review prior to submission to Earth Tech for data validation.

During the validation process, laboratory-qualified and unqualified data are verified against all available supporting documentation. Based on this evaluation, qualifier codes may be added, deleted or modified by the data user. Final results are therefore, either qualified or unqualified. Validator-qualified results are annotated with the following codes in accordance with the Functional Guidelines:

- U The constituent was analyzed for, but was not detected above the level of the associated analytical reporting limit. The associated value is either the sample quantitation limit or the sample detection limit.
- J The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample.
- UJ The analyte was not detected above the reported sample quantitation limit. However, the reported quantitation limit is approximate and may or may not represent the actual limit of quantitation necessary to accurately and precisely measure the analyte in the sample.

- R The sample results are rejected due to serious deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte cannot be verified.

Details of the pesticide data validation findings and conclusions are provided in the following sections of this report:

1. Holding Times
2. Gas Chromatograph/Electronic Capture Detector (GC/ECD) Instrument Performance Check
3. IC
4. Calibration Verification
5. Blanks
6. Surrogate Spikes
7. Matrix Spike/Matrix Spike Duplicate (MS/MSD)
8. Pesticide Cleanup Checks
9. Target Compound Identification
10. Constituent Quantitation and Reported Detection Limits
11. Documentation
12. Overall Assessment

1. HOLDING TIMES

All samples were extracted within the seven-day technical holding time and the five-day VTSR method holding time. The cooler temperature upon receipt at the laboratory was within the recommended temperature of 4°C +/- 2°C.

2. GC/ECD INSTRUMENT PERFORMANCE CHECK

The Performance Evaluation Mixture (PEM) was analyzed at the correct frequency. Absolute retention times were within limits. The percent resolution between adjacent peaks was within QC limits for the Pesticide Analyte Resolution Check. The percent resolution between adjacent peaks is within QC limits for the Performance Evaluation Mixtures (PEM).

The percent breakdown for both 4,4-DDT and Endrin in each PEM was less than 20.0% for both GC columns. The combined percent breakdown for 4,4-DDT and Endrin in each PEM was less than 30.0% for both GC columns.

3. INITIAL CALIBRATION

Individual standard mixtures A and B were analyzed at the correct frequencies and concentrations. The percent resolution criterion was met for Individual standard mixtures A and B.

The Percent Relative Standard Deviation (%RSD) of the calibration factors for each of the single component pesticides was less than 20%.

The multi-component target compounds were analyzed separately on both columns at a single concentration level. Retention times were determined from a minimum of three peaks.

4. CALIBRATION VERIFICATION

Absolute retention times were within appropriate time retention windows

5. BLANKS

One laboratory method blank and equipment blank were analyzed with this SDG. The results are summarized below.

Method Blank 151585

No constituents were detected above the laboratory-reporting limit. This blank corresponds to all samples extracted on 3/8/04.

Equipment Blank SKSWDEB1009

No constituents were detected above the laboratory-reporting limit in the equipment blank collected on 3/2/04.

6. SURROGATE SPIKES

Decachlorobiphenyl (DCB) and tetrachloro-m-xylene (TCX) surrogate spike recoveries were within the acceptance criteria for all samples except for TCX and DCB associated with sample SKSWD031009. There were no target compounds detected in sample SKSWD031009 therefore no action was taken.

7. MATRIX SPIKE/MATRIX SPIKE DUPLICATE

There were no samples submitted for MS/MSD analysis during this sampling event.

8. PESTICIDE CLEANUP CHECKS

Recoveries of all pesticides and surrogates were within 80-120% for the lot of Florisil cartridges utilized for pesticide cleanup with the exception of Endrin (135%), 4,4'-DDT (139%) and Methoxychlor (121%). There were no target compounds detected in the associated samples therefore no action was taken.

9. TARGET COMPOUND IDENTIFICATION

All reported pesticide data were correctly identified with supporting chromatograms present in the data package.

10. CONSTITUENT QUANTITATION AND REPORTED DETECTION LIMITS

Constituent quantitations were correctly calculated and reported for pesticide constituents.

11. DOCUMENTATION

The documentation appeared accurate and in order.

12. OVERALL ASSESSMENT

The results are acceptable with the validator-added qualifiers.

REFERENCES

US EPA, 1994. *National Functional Guidelines for Inorganic Data Review.*

US EPA, 1999. *National Functional Guidelines for Organic Data Review.*

ANALYTICAL RESULTS

PERFORMED BY

GULF COAST ANALYTICAL LABORATORIES, INC.

Report Date 03/31/2004

GCAL Report 204030804

ADDENDUM

Deliver To Earth Tech
200 Vine Street
Wilder, KY 41076
859-442-2300

Attn Pat Higgins

Customer Earth Tech

Project Skinner Landfill

CASE NARRATIVE

Client: Earth Tech **Report:** 204030804

Gulf Coast Analytical Laboratories received and analyzed the sample(s) listed on the sample cross-reference page of this report. Receipt of the sample(s) is documented by the attached chain of custody. This applies only to the sample(s) listed in this report. No sample integrity or quality control exceptions were identified unless noted below.

The ILM04.1 – CLP analysis is resubmitted as an addendum to include an expanded list of compounds at the request of the client. The Addendum includes an updated case narrative.

SEMI-VOLATILES MASS SPECTROMETRY

In the prep batch 270522, the MS/MSD exhibited sporadic recovery failures.

SEMI-VOLATILES GAS CHROMATOGRAPHY

In the Pesticide Florisil check analysis, the recoveries for Endrin, 4,4-DDT and Methoxychor were above recovery limits; however, these compounds were not detected in the associated samples.

In the Pesticide analysis for sample 20403080401 (SKSWD031009), the surrogate Tetrachloro-m-xylene and Decachlorobiphenyl are above the control limits.

METALS

In the ILM04.1 - CLP Metals analysis for prep batches 270882 and 270883, the MS recovery was outside the control limits for Selenium. The LCS recovery was within control limits. This indicates the analysis is in control and the sample is affected by matrix interference. A post-digestion spike was performed on the QC sample for these batches with a recovery of 88% and 99%.

Zinc is flagged as estimated for samples associated with prep batch 270882 due to the fact that the percent difference between the original sample result and the serial dilution result is greater than 10. A chemical or physical interference is suspected.

Aluminum and Manganese is flagged as estimated for samples associated with prep batch 270883 due to the fact that the percent difference between the original sample result and the serial dilution result is greater than 10. A chemical or physical interference is suspected.

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RESUBMITTED

U.S. EPA - CLP
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INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKSWD031009

Lab Name: PROJ AAH GCAL Contract: _____
 Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: _____
 Matrix: (soil / water) Water Lab Sample ID: 20403080401
 Level: (low / med) _____ Date Received: 03/06/04
 % Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum | 1800 | | E | P |
| 7440-36-0 | Antimony | 5.2 | B | | P |
| 7440-38-2 | Arsenic | 2.9 | U | | P |
| 7440-39-3 | Barium | 40.0 | B | | P |
| 7440-41-7 | Beryllium | 0.2 | B | | P |
| 7440-43-9 | Cadmium | 0.2 | U | | P |
| 7440-70-2 | Calcium | 131000 | | | P |
| 7440-47-3 | Chromium | 1.4 | B | | P |
| 7440-48-4 | Cobalt | 1.5 | B | | P |
| 7440-50-8 | Copper | 11.0 | B | | P |
| 7439-89-6 | Iron | 2200 | | | P |
| 7439-92-1 | Lead | 1.5 | U | | P |
| 7439-95-4 | Magnesium | 31600 | | | P |
| 7439-96-5 | Manganese | 87.5 | | | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 2.5 | B | | P |
| 7440-09-7 | Potassium | 4170 | B | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-23-5 | Sodium | 12600 | | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-62-2 | Vanadium | 2.2 | B | | P |
| 7440-66-6 | Zinc | 14.6 | B | E | P |
| 57-12-5 | Cyanide | 0.8 | B | | AS |

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Color Before: COLORLESS Clarity Before: CLEAR Texture: _____
 Color After: COLORLESS Clarity After: CLEAR Artifacts: _____
 Comments: _____

U.S. EPA - CLP
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INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKSWD03D1009

Lab Name: PROJ AAH GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: _____
Matrix: (soil / water) Water Lab Sample ID: 20403080402
Level: (low / med) _____ Date Received: 03/06/04
% Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum | 1750 | | E | P |
| 7440-36-0 | Antimony | 3.7 | U | | P |
| 7440-38-2 | Arsenic | 5.3 | B | | P |
| 7440-39-3 | Barium | 39.1 | B | | P |
| 7440-41-7 | Beryllium | 0.2 | B | | P |
| 7440-43-9 | Cadmium | 0.2 | U | | P |
| 7440-70-2 | Calcium | 124000 | | | P |
| 7440-47-3 | Chromium | 1.3 | B | | P |
| 7440-48-4 | Cobalt | 1.2 | B | | P |
| 7440-50-8 | Copper | 4.5 | B | | P |
| 7439-89-6 | Iron | 2220 | | | P |
| 7439-92-1 | Lead | 1.5 | U | | P |
| 7439-95-4 | Magnesium | 30200 | | | P |
| 7439-96-5 | Manganese | 90.9 | | | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 2.8 | B | | P |
| 7440-09-7 | Potassium | 3960 | B | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-23-5 | Sodium | 12200 | | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-62-2 | Vanadium | 2.2 | B | | P |
| 7440-66-6 | Zinc | 7.2 | B | E | P |
| 57-12-5 | Cyanide | 2.5 | B | | AS |

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Color After: COLORLESS Clarity After: CLEAR Artifacts: _____
Comments: _____

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INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKSWD03MS1009

Lab Name: PROJ AAH GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: _____
Matrix: (soil / water) Water Lab Sample ID: 20403080403
Level: (low / med) _____ Date Received: 03/06/04
% Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum | 3820 | | E | P |
| 7440-36-0 | Antimony | 109 | | | P |
| 7440-38-2 | Arsenic | 48.3 | | | P |
| 7440-39-3 | Barium | 1980 | | | P |
| 7440-41-7 | Beryllium | 49.6 | | | P |
| 7440-43-9 | Cadmium | 48.0 | | | P |
| 7440-70-2 | Calcium | 125000 | | | P |
| 7440-47-3 | Chromium | 197 | | | P |
| 7440-48-4 | Cobalt | 461 | | | P |
| 7440-50-8 | Copper | 247 | | | P |
| 7439-89-6 | Iron | 3090 | | | P |
| 7439-92-1 | Lead | 18.3 | | | P |
| 7439-95-4 | Magnesium | 30700 | | | P |
| 7439-96-5 | Manganese | 573 | | | P |
| 7439-97-6 | Mercury | 4.9 | | | AV |
| 7440-02-0 | Nickel | 464 | | | P |
| 7440-09-7 | Potassium | 4040 | B | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 47.7 | | | P |
| 7440-23-5 | Sodium | 12100 | | | P |
| 7440-28-0 | Thallium | 43.2 | | | P |
| 7440-62-2 | Vanadium | 499 | | | P |
| 7440-66-6 | Zinc | 452 | | E | P |
| 57-12-5 | Cyanide | 90.9 | | | AS |

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Color After: COLORLESS Clarity After: CLEAR Artifacts: _____
Comments: _____

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INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKSWD03DUP1009

Lab Name: PROJ AAH GCAL Contract: _____
 Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: _____
 Matrix: (soil / water) Water Lab Sample ID: 20403080405
 Level: (low / med) _____ Date Received: 03/06/04
 % Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum | 1940 | | E | P |
| 7440-36-0 | Antimony | 3.7 | U | | P |
| 7440-38-2 | Arsenic | 2.9 | U | | P |
| 7440-39-3 | Barium | 40.8 | B | | P |
| 7440-41-7 | Beryllium | 0.2 | B | | P |
| 7440-43-9 | Cadmium | 0.2 | U | | P |
| 7440-70-2 | Calcium | 127000 | | | P |
| 7440-47-3 | Chromium | 1.6 | B | | P |
| 7440-48-4 | Cobalt | 1.4 | B | | P |
| 7440-50-8 | Copper | 8.4 | B | | P |
| 7439-89-6 | Iron | 2170 | | | P |
| 7439-92-1 | Lead | 1.5 | U | | P |
| 7439-95-4 | Magnesium | 31000 | | | P |
| 7439-96-5 | Manganese | 86.7 | | | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 2.8 | B | | P |
| 7440-09-7 | Potassium | 4130 | B | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-23-5 | Sodium | 12200 | | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-62-2 | Vanadium | 2.4 | B | | P |
| 7440-66-6 | Zinc | 12.5 | B | E | P |
| 57-12-5 | Cyanide | 0.5 | U | | AS |

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 Color After: COLORLESS Clarity After: CLEAR Artifacts: _____
 Comments: _____

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INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKSWDEB1009

Lab Name: PROJ AAH GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: _____
Matrix: (soil / water) Water Lab Sample ID: 20403080406
Level: (low / med) _____ Date Received: 03/06/04
% Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum | 25.8 | U | E | P |
| 7440-36-0 | Antimony | 3.7 | U | | P |
| 7440-38-2 | Arsenic | 2.9 | U | | P |
| 7440-39-3 | Barium | 0.3 | U | | P |
| 7440-41-7 | Beryllium | 0.1 | U | | P |
| 7440-43-9 | Cadmium | 0.2 | U | | P |
| 7440-70-2 | Calcium | 27.1 | B | | P |
| 7440-47-3 | Chromium | 0.8 | U | | P |
| 7440-48-4 | Cobalt | 0.4 | U | | P |
| 7440-50-8 | Copper | 3.5 | B | | P |
| 7439-89-6 | Iron | 14.1 | U | | P |
| 7439-92-1 | Lead | 1.5 | U | | P |
| 7439-95-4 | Magnesium | 36.7 | U | | P |
| 7439-96-5 | Manganese | 0.2 | U | | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 0.7 | U | | P |
| 7440-09-7 | Potassium | 42.1 | U | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-23-5 | Sodium | 85.3 | B | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-62-2 | Vanadium | 0.8 | U | | P |
| 7440-66-6 | Zinc | 3.7 | B | E | P |
| 57-12-5 | Cyanide | 1.9 | B | | AS |

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Color After: COLORLESS Clarity After: CLEAR Artifacts: _____
Comments:

U.S. EPA - CLP
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INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKSWD031009 (DISS)

Lab Name: PROJ AAH GCAL Contract: _____
 Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: _____
 Matrix: (soil / water) Water Lab Sample ID: 20403080408
 Level: (low / med) _____ Date Received: 03/06/04
 % Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum | 34.6 | B | | P |
| 7440-36-0 | Antimony | 3.7 | U | | P |
| 7440-38-2 | Arsenic | 5.3 | B | | P |
| 7440-39-3 | Barium | 29.8 | B | | P |
| 7440-41-7 | Beryllium | 0.2 | B | | P |
| 7440-43-9 | Cadmium | 0.2 | U | | P |
| 7440-70-2 | Calcium | 125000 | | | P |
| 7440-47-3 | Chromium | 0.8 | U | | P |
| 7440-48-4 | Cobalt | 0.4 | U | | P |
| 7440-50-8 | Copper | 4.6 | B | | P |
| 7439-89-6 | Iron | 17.2 | B | | P |
| 7439-92-1 | Lead | 1.5 | U | | P |
| 7439-95-4 | Magnesium | 30400 | | | P |
| 7439-96-5 | Manganese | 3.0 | B | | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 1.4 | B | | P |
| 7440-09-7 | Potassium | 3570 | B | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-23-5 | Sodium | 12200 | | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-62-2 | Vanadium | 0.8 | U | | P |
| 7440-66-6 | Zinc | 0.6 | U | | P |

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 Comments: _____

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INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKSWD03D1009 (DISS)

Lab Name: PROJ AAH GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: _____
Matrx: (soil / water) Water Lab Sample ID: 20403080409
Level: (low / med) _____ Date Received: 03/06/04
% Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum | 25.8 | U | | P |
| 7440-36-0 | Antimony | 3.7 | U | | P |
| 7440-38-2 | Arsenic | 4.1 | B | | P |
| 7440-39-3 | Barium | 29.8 | B | | P |
| 7440-41-7 | Beryllium | 0.1 | B | | P |
| 7440-43-9 | Cadmium | 0.2 | U | | P |
| 7440-70-2 | Calcium | 125000 | | | P |
| 7440-47-3 | Chromium | 0.8 | U | | P |
| 7440-48-4 | Cobalt | 0.5 | B | | P |
| 7440-50-8 | Copper | 1.2 | U | | P |
| 7439-89-6 | Iron | 27.7 | B | | P |
| 7439-92-1 | Lead | 1.5 | U | | P |
| 7439-95-4 | Magnesium | 30500 | | | P |
| 7439-96-5 | Manganese | 3.0 | B | | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 1.3 | B | | P |
| 7440-09-7 | Potassium | 3560 | B | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-23-5 | Sodium | 12500 | | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-62-2 | Vanadium | 0.8 | U | | P |
| 7440-66-6 | Zinc | 0.6 | U | | P |

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Color After: COLORLESS Clarity After: CLEAR Artifacts: _____
Comments: _____

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INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKSWD03MS1009(DISS)

Lab Name: PROJ AAH GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: _____
Matrix: (soil / water) Water Lab Sample ID: 20403080410
Level: (low / med) _____ Date Received: 03/06/04
% Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum | 2050 | | | P |
| 7440-36-0 | Antimony | 115 | | | P |
| 7440-38-2 | Arsenic | 48.7 | | | P |
| 7440-39-3 | Barium | 2010 | | | P |
| 7440-41-7 | Beryllium | 50.6 | | | P |
| 7440-43-9 | Cadmium | 49.6 | | | P |
| 7440-70-2 | Calcium | 125000 | | | P |
| 7440-47-3 | Chromium | 201 | | | P |
| 7440-48-4 | Cobalt | 471 | | | P |
| 7440-50-8 | Copper | 247 | | | P |
| 7439-89-6 | Iron | 1020 | | | P |
| 7439-92-1 | Lead | 16.9 | | | P |
| 7439-95-4 | Magnesium | 31000 | | | P |
| 7439-96-5 | Manganese | 500 | | | P |
| 7439-97-6 | Mercury | 5.1 | | | AV |
| 7440-02-0 | Nickel | 475 | | | P |
| 7440-09-7 | Potassium | 3590 | B | | P |
| 7782-49-2 | Selenium | 5.0 | | N | P |
| 7440-22-4 | Silver | 49.3 | | | P |
| 7440-23-5 | Sodium | 12400 | | | P |
| 7440-28-0 | Thallium | 42.3 | | | P |
| 7440-62-2 | Vanadium | 509 | | | P |
| 7440-66-6 | Zinc | 444 | | | P |

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Color After: COLORLESS Clarity After: CLEAR Artifacts: _____
Comments:

U.S. EPA - CLP
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INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKSWD03DUP1009(DISS)

Lab Name: PROJ AAH GCAL Contract: _____
 Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: _____
 Matrix: (soil / water) Water Lab Sample ID: 20403080411
 Level: (low / med) _____ Date Received: 03/06/04
 % Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum | 38.3 | B | | P |
| 7440-36-0 | Antimony | 3.7 | U | | P |
| 7440-38-2 | Arsenic | 2.9 | U | | P |
| 7440-39-3 | Barium | 29.4 | B | | P |
| 7440-41-7 | Beryllium | 0.1 | B | | P |
| 7440-43-9 | Cadmium | 0.2 | U | | P |
| 7440-70-2 | Calcium | 118000 | | | P |
| 7440-47-3 | Chromium | 0.8 | U | | P |
| 7440-48-4 | Cobalt | 1.7 | B | | P |
| 7440-50-8 | Copper | 12.2 | B | | P |
| 7439-89-6 | Iron | 21.5 | B | | P |
| 7439-92-1 | Lead | 1.5 | U | | P |
| 7439-95-4 | Magnesium | 29100 | | | P |
| 7439-96-5 | Manganese | 3.6 | B | | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 3.7 | B | | P |
| 7440-09-7 | Potassium | 3310 | B | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-23-5 | Sodium | 11700 | | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-62-2 | Vanadium | 0.8 | U | | P |
| 7440-66-6 | Zinc | 6.0 | B | | P |

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INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

SKSWDEB1009(DISS)

Lab Name: PROJ AAH GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: _____
Matrix: (soil / water) Water Lab Sample ID: 20403080412
Level: (low / med) _____ Date Received: 03/06/04
% Solids: _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|--------------|----|
| 7429-90-5 | Aluminum | 25.8 | U | | P |
| 7440-36-0 | Antimony | 3.7 | U | | P |
| 7440-38-2 | Arsenic | 2.9 | U | | P |
| 7440-39-3 | Barium | 0.3 | U | | P |
| 7440-41-7 | Beryllium | 0.1 | U | | P |
| 7440-43-9 | Cadmium | 0.2 | U | | P |
| 7440-70-2 | Calcium | 71.9 | B | | P |
| 7440-47-3 | Chromium | 0.8 | U | | P |
| 7440-48-4 | Cobalt | 0.4 | U | | P |
| 7440-50-8 | Copper | 1.2 | U | | P |
| 7439-89-6 | Iron | 22.6 | B | | P |
| 7439-92-1 | Lead | 1.5 | U | | P |
| 7439-95-4 | Magnesium | 36.7 | U | | P |
| 7439-96-5 | Manganese | 0.6 | B | P | P |
| 7439-97-6 | Mercury | 0.1 | U | | AV |
| 7440-02-0 | Nickel | 1.5 | B | | P |
| 7440-09-7 | Potassium | 42.1 | U | | P |
| 7782-49-2 | Selenium | 4.4 | U | N | P |
| 7440-22-4 | Silver | 0.4 | U | | P |
| 7440-23-5 | Sodium | 292 | B | | P |
| 7440-28-0 | Thallium | 2.6 | U | | P |
| 7440-62-2 | Vanadium | 0.8 | U | | P |
| 7440-66-6 | Zinc | 3.9 | B | | P |

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9/23/05
MIL

Color Before: COLORLESS Clarity Before: CLEAR Texture: _____
Color After: COLORLESS Clarity After: CLEAR Artifacts: _____
Comments: _____

INITIAL AND CONTINUING CALIBRATION VERIFICATION

Lab Name: GCAL

Contract: _____

Lab Code: LA024 Case No.: _____

SAS No.: _____ SDG No.: 204030804

Calibration Source: 106-61-2 CPI

Instrument ID: ICP5 ICAL ID: 2

Date Analyzed: 03/18/04 Time: 1140

CRDL STANDARD

| Analyte | True | Found | CAL %R | Units | Method | Type |
|-----------|------|-------|--------|-------|----------------------|------|
| Antimony | 120 | 125 | 104 | ug/L | ILM04.1 - CLP Metals | P |
| Arsenic | 20.0 | 23.2 | 116 | ug/L | ILM04.1 - CLP Metals | P |
| Beryllium | 10.0 | 9.00 | 90 | ug/L | ILM04.1 - CLP Metals | P |
| Cadmium | 10.0 | 9.30 | 93 | ug/L | ILM04.1 - CLP Metals | P |
| Chromium | 20.0 | 17.4 | 87 | ug/L | ILM04.1 - CLP Metals | P |
| Cobalt | 100 | 89.8 | 90 | ug/L | ILM04.1 - CLP Metals | P |
| Copper | 50.0 | 31.5 | 63 | ug/L | ILM04.1 - CLP Metals | P |
| Lead | 6.00 | 6.10 | 101 | ug/L | ILM04.1 - CLP Metals | P |
| Manganese | 30.0 | 28.0 | 93 | ug/L | ILM04.1 - CLP Metals | P |
| Nickel | 80.0 | 73.3 | 92 | ug/L | ILM04.1 - CLP Metals | P |
| Selenium | 10.0 | 6.20 | 62 | ug/L | ILM04.1 - CLP Metals | P |
| Silver | 20.0 | 18.5 | 92 | ug/L | ILM04.1 - CLP Metals | P |
| Thallium | 20.0 | 20.1 | 100 | ug/L | ILM04.1 - CLP Metals | P |
| Vanadium | 100 | 92.2 | 92 | ug/L | ILM04.1 - CLP Metals | P |
| Zinc | 40.0 | 30.3 | 76 | ug/L | ILM04.1 - CLP Metals | P |

INITIAL AND CONTINUING CALIBRATION VERIFICATION

Lab Name: GCAL Contract: _____
 Lab Ccde: LA024 Case No.: _____ SAS No.: _____ SDG No.: 204030804
 Calibration Source: 106-61-2 CPI Instrument ID: ICP5 ICAL ID: 2
 Date Analyzed: 03/18/04 Time: 1445

CRDL STANDARD

| Analyte | True | Found | CAL %R | Units | Method | Type |
|-----------|------|-------|--------|-------|----------------------|------|
| Antimony | 120 | 127 | 106 | ug/L | ILM04.1 - CLP Metals | P |
| Arsenic | 20.0 | 22.6 | 113 | ug/L | ILM04.1 - CLP Metals | P |
| Beryllium | 10.0 | 8.90 | 89 | ug/L | ILM04.1 - CLP Metals | P |
| Cadmium | 10.0 | 9.40 | 94 | ug/L | ILM04.1 - CLP Metals | P |
| Chromium | 20.0 | 17.8 | 89 | ug/L | ILM04.1 - CLP Metals | P |
| Cobalt | 100 | 89.8 | 90 | ug/L | ILM04.1 - CLP Metals | P |
| Copper | 50.0 | 31.4 | 62.8 | ug/L | ILM04.1 - CLP Metals | P |
| Lead | 6.00 | 4.70 | 78.3 | ug/L | ILM04.1 - CLP Metals | P |
| Manganese | 30.0 | 27.9 | 93 | ug/L | ILM04.1 - CLP Metals | P |
| Nickel | 80.0 | 74.1 | 93 | ug/L | ILM04.1 - CLP Metals | P |
| Selenium | 10.0 | 6.40 | 64 | ug/L | ILM04.1 - CLP Metals | P |
| Silver | 20.0 | 17.5 | 87 | ug/L | ILM04.1 - CLP Metals | P |
| Thallium | 20.0 | 19.7 | 99 | ug/L | ILM04.1 - CLP Metals | P |
| Vanadium | 100 | 91.8 | 92 | ug/L | ILM04.1 - CLP Metals | P |
| Zinc | 40.0 | 30.1 | 75.3 | ug/L | ILM04.1 - CLP Metals | P |

INITIAL AND CONTINUING CALIBRATION VERIFICATION

Lab Name: GCAL

Contract: _____

Lab Code: LA024 Case No.: _____

SAS No.: _____ SDG No.: 204030804

Calibration Source: 106-61-2 CPI

Instrument ID: ICP5 ICAL ID: 2

Date Analyzed: 03/18/04 Time: 1537

CRDL STANDARD

| Analyte | True | Found | CAL %R | Units | Method | Type |
|-----------|------|-------|--------|-------|----------------------|------|
| Antimony | 120 | 130 | 108 | ug/L | ILM04.1 - CLP Metals | P |
| Arsenic | 20.0 | 22.5 | 113 | ug/L | ILM04.1 - CLP Metals | P |
| Beryllium | 10.0 | 9.30 | 93 | ug/L | ILM04.1 - CLP Metals | P |
| Cadmium | 10.0 | 9.90 | 99 | ug/L | ILM04.1 - CLP Metals | P |
| Chromium | 20.0 | 19.6 | 98 | ug/L | ILM04.1 - CLP Metals | P |
| Cobalt | 100 | 94.9 | 95 | ug/L | ILM04.1 - CLP Metals | P |
| Copper | 50.0 | 34.5 | 69 | ug/L | ILM04.1 - CLP Metals | P |
| Lead | 6.00 | 7.40 | 123 | ug/L | ILM04.1 - CLP Metals | P |
| Manganese | 30.0 | 28.9 | 96 | ug/L | ILM04.1 - CLP Metals | P |
| Nickel | 80.0 | 78.0 | 97 | ug/L | ILM04.1 - CLP Metals | P |
| Selenium | 10.0 | 8.40 | 84 | ug/L | ILM04.1 - CLP Metals | P |
| Silver | 20.0 | 19.0 | 95 | ug/L | ILM04.1 - CLP Metals | P |
| Thallium | 20.0 | 23.9 | 120 | ug/L | ILM04.1 - CLP Metals | P |
| Vanadium | 100 | 96.7 | 97 | ug/L | ILM04.1 - CLP Metals | P |
| Zinc | 40.0 | 32.3 | 81 | ug/L | ILM04.1 - CLP Metals | P |

U.S. EPA - CLP

3

BLANKS

Lab Name: PROJ AAH GCAL

Contract: _____

Lab Code: LA024

Case No.: _____

SAS No.: _____

SDG No.: _____

Preparation Blank Matrix: (soil / water) WaterPreparation Blank Concentration Units: (ug/L / mg/kg) ug/L

| Analyte | Initial Calib. Blank (ug/L) | | Continuing Calibration Blank (ug/L) | | | | | | Preparation Blank | | M |
|-----------|-----------------------------|---|-------------------------------------|---|-------|---|------|---|-------------------|---|----|
| | | C | 1 | C | 2 | C | 3 | C | | C | |
| Aluminum | 25.8 | U | 25.8 | U | 25.8 | U | 25.8 | U | 25.800 | U | P |
| Antimony | 11.3 | B | 19.8 | B | 7.5 | B | 20.4 | B | 6.501 | B | P |
| Arsenic | 3.1 | B | 4.5 | B | 4.9 | B | 3.7 | B | 2.900 | U | P |
| Barium | 0.4 | B | 0.3 | B | 0.3 | U | 0.3 | U | 0.300 | U | P |
| Beryllium | 0.1 | U | 0.1 | U | 0.1 | B | 0.1 | U | 0.100 | U | P |
| Cadmium | 0.6 | B | 1.6 | B | 0.3 | B | 3.5 | B | 0.550 | B | P |
| Calcium | 7.5 | U | 7.5 | U | 25.2 | B | 11.7 | B | 7.500 | U | P |
| Chromium | 0.8 | U | 0.8 | U | 0.8 | U | 0.8 | U | 0.800 | U | P |
| Cobalt | 0.4 | U | 0.4 | U | 0.4 | U | 0.7 | B | 0.449 | B | P |
| Copper | 14.0 | B | 21.9 | B | 10.7 | B | 15.8 | B | 9.969 | B | P |
| Iron | 14.1 | U | 17.6 | B | 14.1 | U | 26.8 | B | 14.000 | U | P |
| Lead | 1.5 | U | 2.1 | B | 2.5 | B | 1.5 | U | 1.500 | U | P |
| Magnesium | 36.7 | U | 36.7 | U | 36.7 | U | 36.7 | U | 36.700 | U | P |
| Manganese | 0.4 | B | 0.3 | B | 0.4 | B | 0.3 | B | 0.772 | B | P |
| Mercury | -0.1 | B | -0.1 | B | -0.1 | B | -0.1 | B | 0.100 | U | AV |
| Nickel | 0.7 | U | 1.2 | B | 0.8 | B | 2.2 | B | 0.931 | B | P |
| Potassium | 42.1 | U | 42.1 | U | 42.1 | U | 42.1 | U | 42.100 | U | P |
| Selenium | 4.4 | U | 4.4 | U | 4.4 | U | 4.4 | U | -4.403 | B | P |
| Silver | 0.5 | B | 0.4 | U | -1.3 | B | 0.4 | U | -0.835 | B | P |
| Sodium | 45.4 | U | 45.4 | U | 45.4 | U | 45.4 | U | 85.108 | B | P |
| Thallium | -3.6 | B | 2.6 | U | -2.8 | B | 2.6 | U | -3.681 | B | P |
| Vanadium | 0.8 | U | 0.8 | U | -1.1 | B | -0.9 | B | 0.800 | U | P |
| Zinc | 7.1 | B | 8.9 | B | 7.3 | B | 10.3 | B | 7.264 | B | P |
| Cyanide | 0.500 | U | 0.500 | U | 0.500 | U | | | 0.500 | U | AS |

U.S. EPA - CLP

3

BLANKS

Lab Name: PROJ AAH GCAL

Contract: _____

Lab Code: LA024

Case No.: _____

SAS No.: _____

SDG No.: _____

Preparation Blank Matrix: (soil / water) WaterPreparation Blank Concentration Units: (ug/L / mg/kg) ug/L

| Analyte | Initial Calib. Blank (ug/L) | Continuing Calibration Blank (ug/L) | | | | | | Prepa- ration Blank | C | M |
|-----------|--------------------------------------|-------------------------------------|---|---|---|---|---|---------------------------|---|----|
| | | 4 | C | 5 | C | 6 | C | | | |
| Aluminum | | 25.8 | U | | | | | 25.800 | U | P |
| Antimony | | 19.1 | B | | | | | 3.700 | U | P |
| Arsenic | | 2.9 | U | | | | | 2.990 | B | P |
| Barium | | 0.3 | U | | | | | 0.300 | U | P |
| Beryllium | | 0.1 | U | | | | | 0.100 | U | P |
| Cadmium | | 1.6 | B | | | | | 0.200 | U | P |
| Calcium | | 10.6 | B | | | | | 23.995 | B | P |
| Chromium | | 0.8 | U | | | | | -1.275 | B | P |
| Cobalt | | 1.1 | B | | | | | 0.445 | B | P |
| Copper | | 24.6 | B | | | | | 3.617 | B | P |
| Iron | | 35.9 | B | | | | | 14.000 | U | P |
| Lead | | 2.0 | B | | | | | 1.500 | U | P |
| Magnesium | | 36.7 | U | | | | | 36.700 | U | P |
| Manganese | | 0.2 | B | | | | | 0.332 | B | P |
| Mercury | | | | | | | | 0.100 | U | AV |
| Nickel | | 2.2 | B | | | | | 0.877 | B | P |
| Potassium | | 42.1 | U | | | | | 42.100 | U | P |
| Selenium | | 4.4 | U | | | | | 4.400 | U | P |
| Silver | | 0.4 | U | | | | | -2.198 | B | P |
| Sodium | | 45.4 | U | | | | | 88.654 | B | P |
| Thallium | | -4.0 | B | | | | | -2.923 | B | P |
| Vanadium | | 0.8 | U | | | | | -1.051 | B | P |
| Zinc | | 9.8 | B | | | | | 5.114 | B | P |
| Cyanide | | | | | | | | | | AS |

MS/MSD RECOVERY

Lab Name: PROJ AAH GCAL

Contract: _____

Lab Code: LA024

Case No.: _____

SAS No.: _____

SDG No.: _____

Matrix Spike - EPA Sample No: SKSWD031009SAMPLE NO. : 20403080403

| COMPOUND | UNITS | SPIKE ADDED | SAMPLE CONCENTRATION | MS CONCENTRATION | MS % REC | # | QC. LIMITS |
|-----------|-------|----------------|-------------------------|---------------------|--------------------|---|------------|
| Aluminum | ug/L | 2000 | 1800 | 3820 | 101 | | 75 - 125 |
| Antimony | ug/L | 100 | 5.2 | 109 | 104 | | 75 - 125 |
| Arsenic | ug/L | 40 | 2.9 | 48.3 | 121 114 | | 75 - 125 |
| Barium | ug/L | 2000 | 40 | 1980 | 97 | | 75 - 125 |
| Beryllium | ug/L | 50 | .2 | 49.6 | 99 | | 75 - 125 |
| Cadmium | ug/L | 50 | .2 | 48 | 96 | | 75 - 125 |
| Chromium | ug/L | 200 | 1.4 | 197 | 98 | | 75 - 125 |
| Cobalt | ug/L | 500 | 1.5 | 461 | 92 | | 75 - 125 |
| Copper | ug/L | 250 | 11 | 247 | 94 | | 75 - 125 |
| Iron | ug/L | 1000 | 2200 | 3090 | 89 | | 75 - 125 |
| Lead | ug/L | 20 | 1.5 | 18.3 | 92 | | 75 - 125 |
| Manganese | ug/L | 500 | 87.5 | 573 | 97 | | 75 - 125 |
| Mercury | ug/L | 5 | .1 | 4.9 | 99 | | 75 - 125 |
| Nickel | ug/L | 500 | 2.5 | 464 | 92 | | 75 - 125 |
| Selenium | ug/L | 10 | 4.4 | 4.4 | 0 | N | 75 - 125 |
| Silver | ug/L | 50 | .4 | 47.7 | 95 | | 75 - 125 |
| Thallium | ug/L | 50 | 2.6 | 43.2 | 86 | | 75 - 125 |
| Vanadium | ug/L | 500 | 2.2 | 499 | 99 | | 75 - 125 |
| Zinc | ug/L | 500 | 14.6 | 452 | 88 | | 75 - 125 |
| Cyanide | ug/L | 100 | .8 | 90.9 | 90 | | 75 - 125 |

9/22/08
mcc

Column to be used to flag recovery and RPD values with an asterisk

* Values outside of QC limits

RPD : 0 out of 0 outside limitsSpike Recovery: 1 out of 20 outside limits

MS/MSD RECOVERY

Lab Name: PROJ AAH GCAL

Contract: _____

Lab Code: LA024 Case No.: _____

SAS No.: _____ SDG No.: _____

Matrix Spike - EPA Sample No: SKSWD031009 (DISS)SAMPLE NO. : 20403080410

| COMPOUND | UNITS | SPIKE ADDED | SAMPLE CONCENTRATION | MS CONCENTRATION | MS % REC | # | QC. LIMITS |
|-----------|-------|----------------|-------------------------|---------------------|-------------|---|------------|
| Aluminum | ug/L | 2000 | 34.6 | 2050 | 101 | | 75 - 125 |
| Antimony | ug/L | 100 | 3.7 | 115 | 115 | | 75 - 125 |
| Arsenic | ug/L | 40 | 5.3 | 48.7 | 109 | | 75 - 125 |
| Barium | ug/L | 2000 | 29.8 | 2010 | 99 | | 75 - 125 |
| Beryllium | ug/L | 50 | .2 | 50.6 | 101 | | 75 - 125 |
| Cadmium | ug/L | 50 | .2 | 49.6 | 99 | | 75 - 125 |
| Chromium | ug/L | 200 | .8 | 201 | 101 | | 75 - 125 |
| Cobalt | ug/L | 500 | .4 | 471 | 94 | | 75 - 125 |
| Copper | ug/L | 250 | 4.6 | 247 | 97 | | 75 - 125 |
| Iron | ug/L | 1000 | 17.2 | 1020 | 100 | | 75 - 125 |
| Lead | ug/L | 20 | 1.5 | 16.9 | 84 | | 75 - 125 |
| Manganese | ug/L | 500 | 3 | 500 | 99 | | 75 - 125 |
| Mercury | ug/L | 5 | .1 | 5.1 | 101 | | 75 - 125 |
| Nickel | ug/L | 500 | 1.4 | 475 | 95 | | 75 - 125 |
| Selenium | ug/L | 10 | 4.4 | 5 | 50 | N | 75 - 125 |
| Silver | ug/L | 50 | .4 | 49.3 | 99 | | 75 - 125 |
| Thallium | ug/L | 50 | 2.6 | 42.3 | 85 | | 75 - 125 |
| Vanadium | ug/L | 500 | .8 | 509 | 102 | | 75 - 125 |
| Zinc | ug/L | 500 | .6 | 444 | 89 | | 75 - 125 |

Column to be used to flag recovery and RPD values with an asterisk

* Values outside of QC limits

RPD : 0 out of 0 outside limitsSpike Recovery: 1 out of 19 outside limits

FORM V (PART 1) - IN

000280

RESUBMITTED

U.S. EPA - CLP
5B
POST DIGEST SPIKE SAMPLE RECOVERY

EPA SAMPLE NO.

SKSWD031009PDS

Lab Name: PROJ AAH GCAL

Lab Code: LA024

Case No.: _____

Contract: _____

Matrix: (soil / water) Water

SAS No.: _____

SDG No.: _____

% Solids for Sample: _____

Level: (low / med) _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| Analyte | Control Limit | | Spiked Sample | | Sample | | Spike Added (SA) | % R | Q | M |
|-----------|---------------|--|---------------|---|-------------|---|------------------|-----|---|---|
| | %R | | Result (SSR) | C | Result (SR) | C | | | | |
| Aluminum | | | 6730 | | 1800 | | 5000 | 99 | | P |
| Antimony | | | 475 | | 5.2 | B | 500 | 94 | | P |
| Arsenic | | | 474 | | 2.9 | U | 500 | 95 | | P |
| Barium | | | 492 | | 40 | B | 500 | 90 | | P |
| Beryllium | | | 450 | | .2 | B | 500 | 90 | | P |
| Cadmium | | | 439 | | .2 | U | 500 | 88 | | P |
| Calcium | | | 134000 | | 131000 | | 5000 | 59 | | P |
| Chromium | | | 452 | | 1.4 | B | 500 | 90 | | P |
| Cobalt | | | 430 | | 1.5 | B | 500 | 86 | | P |
| Copper | | | 453 | | 11 | B | 500 | 88 | | P |
| Iron | | | 6590 | | 2200 | | 5000 | 88 | | P |
| Lead | | | 435 | | 1.5 | U | 500 | 87 | | P |
| Magnesium | | | 36000 | | 31600 | | 5000 | 88 | | P |
| Manganese | | | 533 | | 87.5 | | 500 | 89 | | P |
| Nickel | | | 428 | | 2.5 | B | 500 | 85 | | P |
| Potassium | | | 13500 | | 4170 | B | 10000 | 93 | | P |
| Selenium | | | 442 | | 4.4 | U | 500 | 88 | | P |
| Silver | | | 453 | | .4 | U | 500 | 91 | | P |
| Sodium | | | 31000 | | 12600 | | 20000 | 92 | | P |
| Thallium | | | 447 | | 2.6 | U | 500 | 89 | | P |
| Vanadium | | | 460 | | 2.2 | B | 500 | 92 | | P |
| Zinc | | | 425 | | 14.6 | B | 500 | 82 | | P |

Comments:

U.S. EPA - CLP
5B
POST DIGEST SPIKE SAMPLE RECOVERY

EPA SAMPLE NO.

SKSWD031009 (DISS)PDS

Lab Name: PROJ AAH GCAL

Lab Code: LA024

Case No.: _____

Contract: _____

Matrix: (soil / water) Water

SAS No.: _____

SDG No.: _____

% Solids for Sample: _____

Level: (low / med) _____

Concentration Units (ug/L or mg/kg dry weight) : ug/L

| Analyte | Control Limit | | Spiked Sample | | Sample | | Spike Added (SA) | % R | Q | M |
|-----------|---------------|--|---------------|---|-------------|---|------------------|-----|---|---|
| | %R | | Result (SSR) | C | Result (SR) | C | | | | |
| Aluminum | | | 5080 | | 34.6 | B | 5000 | 101 | | P |
| Antimony | | | 511 | | 3.7 | U | 500 | 102 | | P |
| Arsenic | | | 535 | | 5.3 | B | 500 | 106 | | P |
| Barium | | | 536 | | 29.8 | B | 500 | 101 | | P |
| Beryllium | | | 512 | | .2 | B | 500 | 102 | | P |
| Cadmium | | | 493 | | .2 | U | 500 | 99 | | P |
| Calcium | | | 136000 | | 125000 | | 5000 | 220 | | P |
| Chromium | | | 507 | | .8 | U | 500 | 101 | | P |
| Cobalt | | | 483 | | .4 | U | 500 | 97 | | P |
| Copper | | | 511 | | 4.6 | B | 500 | 101 | | P |
| Iron | | | 4990 | | 17.2 | B | 5000 | 99 | | P |
| Lead | | | 489 | | 1.5 | U | 500 | 98 | | P |
| Magnesium | | | 37300 | | 30400 | | 5000 | 136 | | P |
| Manganese | | | 502 | | 3 | B | 500 | 100 | | P |
| Nickel | | | 479 | | 1.4 | B | 500 | 96 | | P |
| Potassium | | | 14100 | | 3570 | B | 10000 | 105 | | P |
| Selenium | | | 497 | | 4.4 | U | 500 | 99 | | P |
| Silver | | | 512 | | .4 | U | 500 | 102 | | P |
| Sodium | | | 33900 | | 12200 | | 20000 | 108 | | P |
| Thallium | | | 496 | | 2.6 | U | 500 | 99 | | P |
| Vanadium | | | 515 | | .8 | U | 500 | 103 | | P |
| Zinc | | | 458 | | .6 | U | 500 | 92 | | P |

Comments:

U.S. EPA - CLP

6

DUPLICATES

EPA SAMPLE NO

SKSWD03DUP1009(DISS)

Lab Name: PROJ AAH GCAL

Lab Code: LA024

Case No.: _____

Contract: _____

Matrix: (soil / water) Water

SAS No.: _____

SDG No.: _____

Level: (low / med) _____

% Solids for Sample: _____

% Solids for Duplicate: _____

Concentration Units (ug/L or mg/kg dry weight) ug/L

| Analyte | Control Limit | Sample (S) | C | Duplicate (D) | C | RPD | Q | M |
|-----------|---------------|------------|---|---------------|---|-----|---|----|
| Aluminum | - | 34.6 | B | 38.3 | B | 10 | | P |
| Antimony | - | 3.7 | U | 3.7 | U | 0 | | P |
| Arsenic | - | 5.3 | B | 2.9 | U | 200 | | P |
| Barium | - | 29.8 | B | 29.4 | B | 1 | | P |
| Beryllium | - | .2 | B | .1 | B | 67 | | P |
| Cadmium | - | .2 | U | .2 | U | 0 | | P |
| Calcium | 0 - 20 | 125000 | | 118000 | | 6 | | P |
| Chromium | - | .8 | U | .8 | U | 0 | | P |
| Cobalt | - | .4 | U | 1.7 | B | 200 | | P |
| Copper | - | 4.6 | B | 12.2 | B | 90 | | P |
| Iron | - | 17.2 | B | 21.5 | B | 22 | | P |
| Lead | - | 1.5 | U | 1.5 | U | 0 | | P |
| Magnesium | 0 - 20 | 30400 | | 29100 | | 4 | | P |
| Manganese | - | 3 | B | 3.6 | B | 18 | | P |
| Mercury | - | .1 | U | .1 | U | 0 | | AV |
| Nickel | - | 1.4 | B | 3.7 | B | 90 | | P |
| Potassium | - | 3570 | B | 3310 | B | 8 | | P |
| Selenium | - | 4.4 | U | 4.4 | U | 0 | | P |
| Silver | - | .4 | U | .4 | U | 0 | | P |
| Sodium | 0 - 5000 | 12200 | | 11700 | | 500 | | P |
| Thallium | - | 2.6 | U | 2.6 | U | 0 | | P |
| Vanadium | - | .8 | U | .8 | U | 0 | | P |
| Zinc | - | .6 | U | 6 | B | 200 | | P |

U.S. EPA - CLP

6

DUPLICATES

EPA SAMPLE NO

SKSWD03DUP1009

Lab Name: PROJ AAH GCAL

Lab Code: LA024

Case No.:

Contract:

Matrix: (soil / water) Water

SAS No.:

SDG No.:

Level: (low / med)

% Solids for Sample:

% Solids for Duplicate:

Concentration Units (ug/L or mg/kg dry weight) ug/L

| Analyte | Control Limit | Sample (S) | C | Duplicate (D) | C | RPD | Q | M |
|-----------|---------------|------------|---|---------------|---|-----|---|----|
| Aluminum | 0 - 20 | 1800 | | 1940 | | 7 | | P |
| Antimony | - | 5.2 | B | 3.7 | U | 200 | | P |
| Arsenic | - | 2.9 | U | 2.9 | U | 0 | | P |
| Barium | - | 40 | B | 40.8 | B | 2 | | P |
| Beryllium | - | .2 | B | .2 | B | 0 | | P |
| Cadmium | - | .2 | U | .2 | U | 0 | | P |
| Calcium | 0 - 20 | 131000 | | 127000 | | 3 | | P |
| Chromium | - | 1.4 | B | 1.6 | B | 13 | | P |
| Cobalt | - | 1.5 | B | 1.4 | B | 7 | | P |
| Copper | - | 11 | B | 8.4 | B | 27 | | P |
| Iron | 0 - 20 | 2200 | | 2170 | | 1 | | P |
| Lead | - | 1.5 | U | 1.5 | U | 0 | | P |
| Magnesium | 0 - 20 | 31600 | | 31000 | | 2 | | P |
| Manganese | 0 - 20 | 87.5 | | 86.7 | | .9 | | P |
| Mercury | - | .1 | U | .1 | U | 0 | | AV |
| Nickel | - | 2.5 | B | 2.8 | B | 11 | | P |
| Potassium | - | 4170 | B | 4130 | B | 1 | | P |
| Selenium | - | 4.4 | U | 4.4 | U | 0 | | P |
| Silver | - | .4 | U | .4 | U | 0 | | P |
| Sodium | 0 - 5000 | 12600 | | 12200 | | 400 | | P |
| Thallium | - | 2.6 | U | 2.6 | U | 0 | | P |
| Vanadium | - | 2.2 | B | 2.4 | B | 9 | | P |
| Zinc | - | 14.6 | B | 12.5 | B | 15 | | P |
| Cyanide | - | .8 | B | .5 | U | 200 | | AS |

FORM VI - IN

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RESUBMITTED

U.S. EPA - CLP
7
LABORATORY CONTROL SAMPLE

Lab Name: PROJ AAH GCAL Contract: _____
Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: _____
Solid LCS Source: _____
Aqueous LCS Source: 310006 HIGH PURITY~323003 HIGH

| Analyte | Aqueous (ug/L) | | | Solid (mg/kg) | | | |
|-----------|----------------|-------|-----|---------------|-------|---|-----|
| | True | Found | % R | True | Found | C | % R |
| Aluminum | 2000 | 2130 | 107 | | | | |
| Antimony | 500 | 564 | 113 | | | | |
| Arsenic | 2000 | 2070 | 103 | | | | |
| Barium | 2000 | 2030 | 101 | | | | |
| Beryllium | 50.0 | 52.0 | 104 | | | | |
| Cadmium | 50.0 | 51.6 | 103 | | | | |
| Calcium | 12500 | 13000 | 104 | | | | |
| Chromium | 200 | 201 | 100 | | | | |
| Cobalt | 500 | 493 | 99 | | | | |
| Copper | 250 | 274 | 110 | | | | |
| Iron | 1000 | 1060 | 106 | | | | |
| Lead | 500 | 509 | 102 | | | | |
| Magnesium | 12500 | 13000 | 104 | | | | |
| Manganese | 500 | 508 | 102 | | | | |
| Nickel | 500 | 501 | 100 | | | | |
| Potassium | 12500 | 12400 | 99 | | | | |
| Selenium | 2000 | 2000 | 100 | | | | |
| Silver | 50.0 | 50.4 | 101 | | | | |
| Sodium | 12500 | 12800 | 102 | | | | |
| Thallium | 2000 | 2020 | 101 | | | | |
| Vanadium | 500 | 519 | 104 | | | | |
| Zinc | 500 | 469 | 94 | | | | |

U.S. EPA - CLP
7
LABORATORY CONTROL SAMPLE

Lab Name: PROJ AAH GCAL Contract: _____
 Lab Code: LA024 Case No.: _____ SAS No.: _____ SDG No.: _____

Solid LCS Source: _____

Aqueous LCS Source: 310006 HIGH PURITY~323003 HIGH

| Analyte | Aqueous (ug/L) | | | Solid (mg/kg) | | | |
|-----------|----------------|-------|-----|---------------|-------|---|-----|
| | True | Found | % R | True | Found | C | % R |
| Aluminum | 2000 | 2200 | 110 | | | | |
| Antimony | 500 | 587 | 117 | | | | |
| Arsenic | 2000 | 2160 | 108 | | | | |
| Barium | 2000 | 2100 | 105 | | | | |
| Beryllium | 50.0 | 53.9 | 108 | | | | |
| Cadmium | 50.0 | 54.0 | 108 | | | | |
| Calcium | 12500 | 13600 | 109 | | | | |
| Chromium | 200 | 210 | 105 | | | | |
| Cobalt | 500 | 511 | 102 | | | | |
| Copper | 250 | 288 | 115 | | | | |
| Iron | 1000 | 1080 | 108 | | | | |
| Lead | 500 | 530 | 106 | | | | |
| Magnesium | 12500 | 13500 | 108 | | | | |
| Manganese | 500 | 526 | 105 | | | | |
| Nickel | 500 | 520 | 104 | | | | |
| Potassium | 12500 | 13000 | 104 | | | | |
| Selenium | 2000 | 2080 | 104 | | | | |
| Silver | 50.0 | 54.3 | 109 | | | | |
| Sodium | 12500 | 13500 | 108 | | | | |
| Thallium | 2000 | 2110 | 105 | | | | |
| Vanadium | 500 | 538 | 108 | | | | |
| Zinc | 500 | 491 | 98 | | | | |

U.S. EPA - CLP
9
ICP SERIAL DILUTIONS

EPA SAMPLE NO.

SKSWD031009SD

Lab Name: PROJ AAH GCAL

Lab Code: LA024

Case No.

Contract:

Matrix: (soil / water) Water

SAS No.:

SDG No.:

Level: (low / med)

Concentration Units: ug/L

| Analyte | Initial Sample Result (I) | C | Serial Dilution Result (S) | C | % Difference | Q | M |
|-----------|---------------------------------|---|----------------------------------|---|--------------|---|---|
| Aluminum | 1800 | | 2330 | | 29.4 | E | P |
| Antimony | 5.2 | B | 22.6 | B | 335 | | P |
| Arsenic | 2.9 | U | 21.3 | B | | | P |
| Barium | 40.0 | B | 41.9 | B | 4.8 | | P |
| Beryllium | 0.2 | B | 0.5 | U | 150 | | P |
| Cadmium | 0.2 | U | 1.0 | U | | | P |
| Calcium | 131000 | | 132000 | | .8 | | P |
| Chromium | 1.4 | B | 4.0 | U | 186 | | P |
| Cobalt | 1.5 | B | 2.0 | U | 33.3 | | P |
| Copper | 11.0 | B | 6.0 | U | 45.5 | | P |
| Iron | 2200 | | 2370 | | 7.7 | | P |
| Lead | 1.5 | U | 7.5 | U | | | P |
| Magnesium | 31600 | | 31700 | | .3 | | P |
| Manganese | 87.5 | | 91.7 | | 4.8 | | P |
| Nickel | 2.5 | B | 3.6 | B | 44 | | P |
| Potassium | 4170 | B | 3990 | B | 4.3 | | P |
| Selenium | 4.4 | U | 22.0 | U | | | P |
| Silver | 0.4 | U | 2.0 | U | | | P |
| Sodium | 12600 | | 12100 | B | 4 | | P |
| Thallium | 2.6 | U | 13.0 | U | | | P |
| Vanadium | 2.2 | B | 4.0 | U | 81.8 | | P |
| Zinc | 14.6 | B | 3.0 | U | 79.5 | E | P |

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9
ICP SERIAL DILUTIONS

EPA SAMPLE NO.

SKSWD031009 (DISS)SD

Lab Name: PROJ AAH GCAL

Lab Code: LA024

Case No.

Contract:

Matrix: (soil / water) Water

SAS No.:

SDG No.:

Level: (low / med)

Concentration Units: ug/L

| Analyte | Initial Sample Result (I) | C | Serial Dilution Result (S) | C | % Difference | Q | M |
|-----------|---------------------------------|---|----------------------------------|---|--------------|---|---|
| Aluminum | 34.6 | B | 148 | B | 328 | | P |
| Antimony | 3.7 | U | 18.5 | U | | | P |
| Arsenic | 5.3 | B | 14.5 | U | 174 | | P |
| Barium | 29.8 | B | 28.5 | B | 4.4 | | P |
| Beryllium | 0.2 | B | 0.5 | U | 150 | | P |
| Cadmium | 0.2 | U | 1.0 | U | | | P |
| Calcium | 125000 | | 126000 | | .8 | | P |
| Chromium | 0.8 | U | 4.0 | U | | | P |
| Cobalt | 0.4 | U | 2.0 | U | | | P |
| Copper | 4.6 | B | 6.0 | U | 30.4 | | P |
| Iron | 17.2 | B | 86.8 | B | 405 | | P |
| Lead | 1.5 | U | 7.5 | U | | | P |
| Magnesium | 30400 | | 30500 | | .3 | | P |
| Manganese | 3.0 | B | 7.2 | B | 140 | | P |
| Nickel | 1.4 | B | 3.5 | U | 150 | | P |
| Potassium | 3570 | B | 3290 | B | 7.8 | | P |
| Selenium | 4.4 | U | 22.0 | U | | | P |
| Silver | 0.4 | U | 2.0 | U | | | P |
| Sodium | 12200 | | 12000 | B | 1.6 | | P |
| Thallium | 2.6 | U | 13.0 | U | | | P |
| Vanadium | 0.8 | U | 4.0 | U | | | P |
| Zinc | 0.6 | U | 3.0 | U | | | P |

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INSTRUMENT DETECTION LIMITS (QUARTERLY)

Lab Name: PROJ AAH GCALLab Code: LA024

Case No.: _____

SAS No.: _____

SDG No.: _____

Instrument ID: ICP5Study Date: 04/30/03

| <i>Analyte</i> | <i>Wavelength (nm)</i> | <i>Background</i> | <i>CRDL (ug/L)</i> | <i>IDL (ug/L)</i> | <i>M</i> |
|----------------|----------------------------|-------------------|------------------------|-----------------------|----------|
| Aluminum | 308.210 | | 200 | 25.8 | P |
| Antimony | 206.830 | | 60 | 3.7 | P |
| Arsenic | 193.700 | | 10 | 2.9 | P |
| Barium | 233.520 | | 200 | .3 | P |
| Beryllium | 313.100 | | 5 | .1 | P |
| Cadmium | 214.430 | | 5 | .2 | P |
| Calcium | 315.880 | | 5000 | 7.5 | P |
| Chromium | 267.710 | | 10 | .8 | P |
| Cobalt | 228.610 | | 50 | .4 | P |
| Copper | 324.750 | | 25 | 1.2 | P |
| Iron | 259.940 | | 100 | 14.1 | P |
| Lead | 220.350 | | 3 | 1.5 | P |
| Magnesium | 279.080 | | 5000 | 36.7 | P |
| Manganese | 257.610 | | 15 | .2 | P |
| Nickel | 231.600 | | 40 | .7 | P |
| Potassium | 766.480 | | 5000 | 42.1 | P |
| Selenium | 196.030 | | 5 | 4.4 | P |
| Silver | 328.060 | | 10 | .4 | P |
| Sodium | 589.580 | | 5000 | 45.4 | P |
| Thallium | 190.800 | | 10 | 2.6 | P |
| Vanadium | 290.880 | | 50 | .8 | P |
| Zinc | 213.860 | | 20 | .6 | P |

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ANALYSIS RUN LOG

Lab Name: PROJ AAH GCAL

Contract: _____

Lab Code: LA024 Case No.: _____

SAS No.: _____ SDG No.: _____

Instrument ID Number: ICP5

Method Type: P

Start Date: 03/18/04

End Date: 03/18/04

Analyte Symbols

| EPA Sample No. | D/F | Time | % R | Al | Sb | As | Ba | Be | Cd | Ca | Cr | Co | Cu | Fe | Pb | Mg | Mn | Hg | Ni | K | Se | Ag | Na | Tl | V | Zn | Cn |
|----------------------|-----|------|-----|----|----|----|----|----|----|----|----|----|----|----|----|----|----|----|----|---|----|----|----|----|---|----|----|
| ICV | 1 | 1108 | | | X | X | | X | X | X | X | X | X | X | X | X | X | | X | | X | X | | X | X | X | |
| ICV2 | 1 | 1114 | | X | | | X | | | | | | | | | | | | | X | | | X | | | | |
| ICB | 1 | 1133 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| CRDL | 1 | 1140 | | | X | X | | X | X | | X | X | X | | X | | X | | X | | X | X | | X | X | X | |
| ICSA | 1 | 1200 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| ICSAB | 1 | 1206 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| CCV | 1 | 1212 | | X | | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| CCV2 | 1 | 1218 | | | X | | | | | | | | | | | | | | | | | | | | | | |
| CCB | 1 | 1225 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| MB153201 | 1 | 1232 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKSWD031009 | 1 | 1239 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKSWD03DUP1009 | 1 | 1245 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKSWD031009SD | 5 | 1252 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKSWDEB1009 | 1 | 1259 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKSWD03D1009 | 1 | 1306 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKSWD03MS1009 | 1 | 1312 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKSWD031009PDS | 1 | 1318 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| LCS153202 | 1 | 1324 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| CCV | 1 | 1337 | | X | | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| CCV2 | 1 | 1343 | | | X | | | | | | | | | | | | | | | | | | | | | | |
| CCB | 1 | 1350 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| MB153203 | 1 | 1357 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKSWD031009 (DISS) | 1 | 1404 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKSWD03DUP1009(DISS) | 1 | 1411 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKSWD031009 (DISS)SD | 5 | 1418 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKSWDEB1009(DISS) | 1 | 1425 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKSWD03D1009 (DISS) | 1 | 1432 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |

U.S. EPA - CLP

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ANALYSIS RUN LOG

Lab Name: PROJ AAH GCAL

Contract: _____

Lab Code: LA024 Case No.: _____

SAS No.: _____ SDG No.: _____

Instrument ID Number: ICP5

Method Type: P

Start Date: 03/18/04

End Date: 03/18/04

Analyte Symbols

| EPA Sample No. | D/F | Time | % R | Al | Sb | As | Ba | Be | Cd | Ca | Cr | Co | Cu | Fe | Pb | Mg | Mn | Hg | Ni | K | Se | Ag | Na | Tl | V | Zn | Cn |
|-----------------------|-----|------|-----|----|----|----|----|----|----|----|----|----|----|----|----|----|----|----|----|---|----|----|----|----|---|----|----|
| SKSWD03MS1009(DISS) | 1 | 1439 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| CRDL | 1 | 1445 | | | X | X | | X | X | | X | X | X | | X | | X | | X | | X | X | | X | X | X | |
| ICSA | 1 | 1452 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| ICSAB | 1 | 1458 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| CCV | 1 | 1504 | | X | | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| CCV2 | 1 | 1510 | | | X | | | | | | | | | | | | | | | | | | | | | | |
| CCB | 1 | 1517 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| SKSWD031009 (DISS)PDS | 1 | 1524 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| LCS153204 | 1 | 1531 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| CRDL | 1 | 1537 | | | X | X | | X | X | | X | X | X | | X | | X | | X | | X | X | | X | X | X | |
| ICSA | 1 | 1600 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| ICSAB | 1 | 1606 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| CCV | 1 | 1612 | | X | | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |
| CCV2 | 1 | 1618 | | | X | | | | | | | | | | | | | | | | | | | | | | |
| CCB | 1 | 1625 | | X | X | X | X | X | X | X | X | X | X | X | X | X | X | | X | X | X | X | X | X | X | X | |